## BIOREMEDIATION AND SOIL FLUSHING TREATABILITY STUDY REPORT L. E. CARPENTER AND CO. PART 1 OF 2

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PREPARED BY IT CORPORATION

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## 1.0 Introduction

L.E. Carpenter and Company (Carpenter) requested laboratory treatability testing in support of a Feasibility Study for a Superfund site in Wharton, New Jersey. From 1943 to 1987, the site housed a production facility for the manufacture of vinyl wall coverings. Between 1963 and 1970, polyvinyl chloride sludge was disposed of in a surface impoundment. Additionally, the site contained numerous support facilities including underground storage tanks, an above ground tank farm, various process waste tanks, cooling water equipment, and other processing equipment. The primary contaminants on site are:

- Bis(2-ethyl hexyl)phthalate (DEHP)
- Xylenes
- Lubricating oil
- Ethyl benzene
- Naphtha

Approximately 5,000 gallons of water immiscible material consisting of these compounds have been pumped from the site subsurface since 1984. Primary contaminants in soil and groundwater are phthalates, polynuclear aromatic hydrocarbons, ethyl benzene, and xylenes.

IT Corporation (IT) conducted the described bioremediation and soil flushing treatability studies for the L.E. Carpenter site under a subcontract from Roy F. Weston, Inc. (WESTON).

## 2.0 Remedial Technology Description

#### 2.1 Bioremediation

Bioremediation of the site could employ both in situ and ex situ treatment strategies. In situ bioremediation has been successfully applied to the remediation of contaminated aquifers impacted with hydrocarbons such as benzene, ethyl benzene, lubricating oils, and xylenes. Microorganisms indigenous to the subsurface are most commonly used for the remediation. In practice, mechanisms for the introduction of oxygen and nutrients and control of pH are employed to enhance the subsurface environment to stimulate biological activity. A variety of engineering designs may be employed to ensure proper control of environmental parameters.

Ex situ bioremediation of groundwater would employ a bioreactor. A fixed film reactor was suggested in the Request for Proposal (RFP) as the reactor operating mode. The present scope of work, however, describes batch studies to determine the potential for successful biodegradation of DEHP and other hydrocarbons in groundwater.

#### 2.2 Soil Flushing

Soil flushing is an in situ treatment process that incorporates the infiltration and percolation of water through a soil profile to mobilize chemical constituents within the soil and transport them to areas for accumulation and ultimate collection. The process consists of applying an aqueous solution to an unexcavated soil either through surface applications, such as flooding or sprinklers, or through injection. A series of shallow well points or subsurface drains are then used to collect the flushing solution containing contaminants and return it to the surface for subsequent treatment or disposal. Following treatment, the groundwater can be reinjected, discharged, or applied upgradient of the extraction wells for continued flushing of the soil.

The movement of an aqueous solution through the soil mobilizes chemical constituents through mass action and chemical reaction mechanisms. Aqueous flushing solutions may be amended with selected reagents to enhance chemical solubility, resulting in increased mobility and therefore increased recovery of chemicals in the leachate or groundwater. The addition of surfactants (e.g., anionic, nonionic, cationic, or amphoteric) is one method for amending aqueous flushing solutions used in treating soils contaminated with organic chemicals.

## 3.0 Conclusions and Recommendations

#### 3.1 Initial Characterization

The treatability study determined that bioremediation has the potential to biodegrade the target compounds, ethyl benzene, xylene, and DEHP. Native microbes possess the necessary metabolic capability to degrade ethyl benzene, xylene, and DEHP. The biodegradation of these compounds was demonstrated in bench-scale batch studies; however, biodegradation has not been determined under in situ conditions where complex interactions among water, soil, ions, organic compounds, and microbes exist. The preliminary evaluation of site characteristics indicated that the site is acceptable for in situ biological treatment.

The following conclusions were drawn from the initial analytical characterization of site samples:

- A relatively large microbial population was detected within the impacted aquifer.
- In most cases, microorganisms present in the groundwater and soil responded very positively to nutrient and oxygen stimulation.
- The site soil had moderate to high phosphate levels and moderate to very low ammonium content. Nutrient addition will be required to enhance the biodegradation process as indicated by the nutrient stimulation test.
- Nutrient compatibility was excellent. The recommended nutrient concentration for injection into the aquifer is 10,000 mg of Restore® 375 per liter.
- Nutrient adsorption by site soil was high on the first addition of nutrients; however, subsequent nutrient additions resulted in progressively lower adsorption. Nutrient adsorption should not seriously interfere with nutrient transport after the first few batch additions.
- The efficiency of oxygen delivery using hydrogen peroxide will be acceptable.
- With one exception, the pH of all soil samples was within the preferred range for successful bioremediation. The pH of the groundwater was within the preferred range. Initial adjustment of the pH is not recommended; however, periodic monitoring should be performed to ensure that drastic changes in pH do not occur.

## 3.2 Groundwater Biotreatability Study

The groundwater biotreatability study indicated efficient biological removal of target compounds using indigenous microorganisms. Groundwater batch treatment results are summarized below:

- Greater than 99.9 percent of the ethyl benzene was removed from all biologically active treatments within 43 hours. Greater than 99.9 percent of the xylene was removed from the activated sludge treatment within 43 hours. Xylene removal from the two biologically active groundwater treatments proceeded at a slower rate compared to the sludge-amended treatment with approximately 50 percent removed after 43 hours. Xylene was nondetectable (<6.6 µg/L) in all biologically active treatments after 183 hours. Appreciable loss of ethyl benzene (approx. 80 percent) and xylene (approx. 75 percent) was observed in the abiotic control. This loss was attributed to volatilization.
- DEHP was biodegraded with a half-life of approximately 60 to 70 hours in the treatments
  containing site microbes. DEHP removal efficiency was 99.7 percent. Activated sludge
  was much less efficient in DEHP biodegradation with a calculated DEHP half-life of 176
  hours and a removal efficiency of 86 percent. Virtually no DEHP was lost from the abiotic
  control.
- No microbial toxicity was observed during the study.
- Microbial enumerations, volatile solids content, and target compound loss generally
  indicate that the addition of microbes to the groundwater does not provide a long-term
  benefit.
- No appreciable changes in nutrient concentration occurred during treatment.
- The pH of groundwater treatments containing native microbes increased slightly whereas the pH of the activated sludge augmented treatment became more acidic during treatment.

#### 3.3 Soil Flushing Treatability Study

The treatability study determined that soil flushing with surfactant-amended flushing solutions has a greater potential to remove DEHP from soil than just the use of potable water. The study also indicated that the type of surfactant may play a vital role in determining the leaching efficiency of DEHP from soil. In this study, a 0.5 percent Brij 30/35 flushing solution was much more effective than a 0.5 percent Tween 85 solution in removing DEHP from soil.

The following conclusions were drawn from the soil flushing treatability study:

- Geotechnical characterization of site soils indicate that the hydraulic conductivity of these soils is high enough to support an in situ remedial process such as soil flushing.
- The plasticity index of these soils indicates that the low cohesive qualities of site soils
  would allow for long term soil flushing with minimum potential for decreases in hydraulic
  conductivity during the remedial process.
- VOCs were readily removed from soil under bench-scale conditions using only potable water and a high percentage of these analytes were removed during the initial stages of soil flushing.
- Phthalates were not responsive to soil flushing using just potable water.

- The use of surfactants to amend the flushing solution increases the amount of phthalates removed in the leachate.
- The type of surfactant used in the flushing solution is a critical factor relative to the efficiency with which phthalates are removed from soil.
- The percent of DEHP removed from soil can be relatively high, even when high concentrations of the compound are initially present within the soil.

#### 3.4 Recommendations

The results of laboratory treatability studies indicate that both soil flushing and bioremediation have potential benefits for site remediation. In situ aquifer bioremediation is technically feasible and should continue to be considered as a candidate for treatment of this site. Hydraulic control of the contaminant plume and the development of a groundwater injection and recovery system that will permit treatment of the plume is critical for successful in situ aquifer bioremediation. An above ground bioreactor may be a viable technology to remove residual organics from recovered groundwater. Site microbes were shown to be effective in biodegrading ethyl benzene, xylene, and DEHP.

Soil flushing with water was effective in removing VOCs; however, water alone was ineffective for flushing DEHP. A surfactant-amended flushing solution was effective in removing DEHP from soil. Soil flushing, particularly with surfactant amended solutions, may be an effective technology for reducing the concentration of VOCs and DEHP in the soil.

A combination of soil flushing and bioremediation may provide an efficient treatment strategy for remediating the subsurface. Surfactant usage must be carefully controlled because many surfactants are biodegradable and, thus, they represent a significant oxygen demand. They may also be preferentially biodegraded leading to reduced efficiency in target compound removal.

## 4.0 Treatability Study Approach

## 4.1 Test Objectives and Rationale

The treatability study evaluated the potential applicability of bioremediation and soil flushing as remedial technologies for contaminant removal from groundwater and soils at the Carpenter site. The primary focus of the remedy screening tests was to evaluate the efficacy of these two technologies on the removal of designated target compounds. Chemical constituents targeted for monitoring throughout these remedy screening studies were DEHP, xylene, and ethyl benzene. The amount of each targeted chemical constituent removed, relative to the initial amounts in groundwater and soil, was used to evaluate the effectiveness of the technologies. The main goals of the study were (1) to show the effectiveness of the treatment strategies in reducing the concentrations of target compounds, and (2) to provide design information required for the next level of testing.

## 4.2 Experimental Design and Procedures

The experimental design provided data for evaluating changing concentrations of target compounds, microbes, nutrients, and pH (Tables 1 and 2). Each treatment scheme was run in triplicate to permit statistical comparison of each treatment's performance.

For respirometric studies of compound biodegradation, oxygen consumption was monitored for the duration of the study. The total mass of carbon dioxide produced during the study was determined at the end of the study by measuring the inorganic carbon content of a potassium hydroxide (KOH) trapping solution in each respirometer vessel.

Biologically-inhibited controls treated with mercuric chloride were used to determine compound loss due to physical processes such as volatilization and adsorption to the reaction vessel. A zero headspace untreated control was established to quantify the loss of target compounds and residual nutrients and the change in microbial density and pH in a completely untreated control. The experimental design accommodates conservation of the mass balance within each treatment vessel and provides for evaluating losses by biological and physical mechanisms.

Reduction in target compound concentration at the end of the study was evaluated using single classification Model I analysis of variance (ANOVA) (Sokal, R. R. and F. J. Rohlf, 1981, Biometry: The Principles and Practice of Statistics in Biological Research, 2nd Edition, W. H. Freeman and Co., San Francisco). The mean residual concentrations of ethyl benzene, xylene, and DEHP were compared for each treatment to determine if a significant difference at the 95 percent confidence level could be attributed to treatment. The null hypothesis states that there is no difference among means of the various treatments. Additionally, planned comparisons were

conducted using the F-test to determine if significant differences at the 95% confidence limit existed between treatments.

#### 4.3 Equipment, Materials, and Methods

All treatments to be evaluated for target compound biodegradation or physical removal were conducted in clean glass vessels or Teflon-lined containers. Abiotic controls were established and treated exactly like biologically active controls to quantify abiotic loss of target compounds.

#### 4.3.1 Biotreatability Equipment and Materials

DEHP, xylenes, and ethyl benzene were analyzed using gas chromatography (GC). Methods were developed for rapid screening of the biotreatability study performance. Split samples, taken at initial and final time points, were submitted to a contract laboratory for analysis by standard procedures, US EPA Method 8240 for volatile organic compounds and US EPA Method 8270 for semivolatiles, to verify initial, intermediate, and final analyses accomplished using rapid turnaround screening procedures.

The total organic carbon (TOC) content of all aqueous treatments was determined using a Dohrmann Total Carbon Analyzer. The pH of treatments was determined using Orion pH electrodes. The nitrogen, as ammonia, and phosphate content of each treatment was determined using modified Standard Method 4500-NH<sub>3</sub> F and 4500-P E, respectively (Standard Methods for the Examination of Water and Wastewater, 17th edition). Microbial respiration was quantified using computerized respirometers (N-Con Systems, Larchmont, New York).

## 4.3.2 Soil Flushing Equipment and Materials

Soil samples were flushed using a flexible wall permeameter and modified control system as described in Section 4.3.2.1.

## 4.3.2.1 Bench-Scale Studies Using a Flexible Wall Permeameter

Bench-scale soil flushing studies are necessary to determine the applicability of a particular site's soil to in situ remediation and to design a soil flushing treatment system. Column studies are the primary method for determining the leachability of chemical constituents from a particular soil. To overcome the adverse effects of a rigid wall column system, a flexible wall permeameter was adapted from geotechnical soil testing and incorporated into a bench-scale soil flushing treatability system.

Undisturbed soil samples collected in the field are used intact in the soil flushing treatability study. Careful consideration is given to establish the flow rate for a specific soil-chemical matrix. Residence time of the flushing solution in the sample is a critical factor in determining the removal efficiency of chemical constituents from the soil. Equilibrium conditions between the chemicals

bound to the soil and their concentrations in the flushing solution are dependent on this residence time as well as solubility and dissolution constants for each chemical. Field-time conditions can be accelerated in the long-term soil flushing study, often to the extent that several decades of fluid flow can be predicted in only a few months of laboratory testing. This task is accomplished by permeating fluids through the sample under higher pressure gradients and greater flow rates than would typically occur in the field environment.

The volume of solution passed through a sample during the study is described in terms of the number of pore volumes flushed through the soil sample per unit of time. The pore volume is easily determined for each sample prior to initiation of flushing studies. Expressing chemical removal on a pore volume basis permits extrapolation of bench-scale data to real-time site estimations for remediation.

## 4.3.2.2 System Design

Figures 1 and 2 illustrate the bench-scale soil flushing treatability test system. The system included a control panel, peristaltic pump, flushing solution feed reservoirs, flexible wall permeameter cells, digital transducer, sample collection reservoirs, resin traps, and a refrigerated leachate collection cabinet. The control panel was used to regulate and monitor cell water and the confining hydraulic pressure exerted on individual soil samples contained within each cell. Separate pressure regulators control the confining cell pressure while cell water volume is monitored by individual burettes. Cell water volume was monitored to ensure that cell water did not enter the sample and contribute to the volume of flushing solution.

The soil flushing solution was supplied to individual cells from separate glass reservoirs. Flow rates were adjusted to deliver approximately five to ten pore volumes per 24-hour flushing period and simultaneously provide sufficient quantites of leachate per collection period for the required analyses. The flushing solution entered the bottom of the cell through the pedestal, was distributed across the width of the sample, and moved in an upward direction through the soil sample. The sample, wrapped in Teflon® and encased in a latex membrane, was confined within the cell by an external pressure applied to the cell water. The flushing solution flowed directly through the sample with minimal movement along the sample/confining wall interface. This flow pattern also minimized any internal channelling of the flushing solution as it moved through the soil.

As the flushing solution started to move up through the sample, air was displaced from the pores and replaced with solution. This resulted in a nearly saturated sample throughout the course of the flushing study. The flushing solution exited the sample through the top cap and flowed via Teflon tubing to the refrigerated leachate collection bottles. Each glass collection bottle was equipped with three Teflon® fittings that allowed leachate entry, leachate overflow to additional collection bottles, and venting of headspace gases through Tenax® resin columns to the atmosphere.

Each permeameter cell was equipped with inlet/outlet monitoring valves in line with a digital pressure transducer. Because flow rate was controlled by a pump, these valves allowed for the head and back pressure for each sample to be monitored constantly throughout the soil flushing study. This feature continuous monitoring of the hydraulic conductivity of each sample during the entire soil flushing study. Each cell was also equipped with sampling ports in line with the inlet and outlet valves. Samples of the flushing solution could be collected as the solution entered and exited the soil sample, at any time during the operation of the flushing study. This feature was used to monitor dissolved oxygen concentration of the soil flushing solution, allowing for an estimation of the level of microbial activity during the flushing study.

#### 4.4 Sampling and Analysis

#### 4.4.1 Experimental Design

Shelby tube samples collected in the field were shipped and stored at 4°C until used in the soil flushing study. Intact soil samples were extruded from the polyethylene liner and trimmed to an approximate size of 7.1-centimeter (cm) diameter by 12-cm length. Representative samples of each soil were analyzed for an initial estimate of the concentration of volatile organic compounds (VOC) and semivolatile organic compounds (SVOC) expected in the actual samples used in the flushing study, as described in Section 4.4.3.4.

The diameter, height, weight, and moisture content were determined for each of the four intact soil samples collected from the site. Specific geotechnical characteristics were calculated from these measurements. Samples were wrapped in Teflon tape and confined in a latex membrane. A filter paper (Whatman No. 42), followed by a porous stone, was attached to each end of the sample. All sample handling was performed in a walk-in cooler at 4°C.

Each sample was placed inside the plexiglass permeameter cell and attached to the pedestal and top cap. All cells were filled with water and a confining pressure was applied to the cell. Each cell was maintained at ambient temperature. The flushing solution was pumped through 1.5 millimeter (mm) PharmMed<sup>TM</sup> tubing to deliver a flow rate of 0.5 milliliter per minute (mL/min) or approximately 5 to 10 pore volumes per 24 hours.

The soil flushing study was maintained for an equivalent 10-year leachate collection period compressed to 60 days. Tenax resin columns were used to collect volatile organics from the headspace of the leachate collection bottles. Leachate samples were collected and stored at 4°C until analyzed. Accumulated leachate was collected for each of the three flushing periods. Three sample analysis periods (TS<sub>0</sub>, TS<sub>1</sub>, and TS<sub>final</sub>) were conducted at the beginning, middle, and end of the soil flushing cycle. Total leachates were analyzed for VOCs (U.S. EPA Method 8240 - gas chromatography/mass spectroscopy [GC/MS]) and SVOCs (U.S. EPA Method 8270 - GC/MS).

Hydraulic conductivity was determined for all samples. The flexible wall permeameter system is designed to allow for monitoring hydraulic conductivity during the course of the soil flushing study. Measurements were taken at specific time intervals throughout the soil flushing treatment in order to monitor fluctuations in hydraulic conductivity during the study.

#### 4.4.2 Collection of Site Samples

## 4.4.2.1 Sample Locations and Collection

Field samples were collected by WESTON using the IT Biotechnology Applications Center (BAC) standard operating procedure for collection and shipping of biological samples. Samples for preliminary analytical screening were submitted directly to IT Analytical Services (ITAS). A report from WESTON describing the locations and depths of the field samples is attached (Appendix A).

#### 4.4.2.2 Groundwater Sampling

Three well locations were selected with respect to the hydrology of the site, an upgradient, downgradient, and source area. These wells have been reported to provide a representative sampling of the full range of VOCs and SVOCs concentrations. VOCs range from nondetectable to  $136,000 \, \mu g/L$ . SVOCs were reported to range from nondetectable to  $67,898 \, \mu g/L$ .

#### 4.4.2.3 Soil Sampling

Three soil sampling locations were selected based on their proximity to previous test pit sampling locations. These sampling locations with documented contaminant concentrations (Remedial Investigation Report) are considered representative of a range in soil contaminant levels (low, medium, and high). Test pits in the area with high levels of contaminants were reported to have concentrations of DEHP and total targeted volatiles ranging from 3,100 to 30,000 mg/kg and 0.2 to 532 mg/kg, respectively.

## 4.4.3 Sampling and Analysis During Treatment

## 4.4.3.1 Groundwater Analyses

Initial Analyses. Groundwater samples provided by Carpenter underwent the following analytical testing:

| Total solids           | U.S. EPA Method 160.3 |
|------------------------|-----------------------|
| Total suspended solids | U.S. EPA Method 160.2 |
| Total dissolved solids | U.S. EPA Method 160.1 |
| Conductivity           | U.S. EPA Method 120.1 |
| Acidity                | U.S. EPA Method 305.1 |
| Alkalinity             | U.S. EPA Method 310.1 |

| рН   | U.S. EPA Method 150.1   |
|--|-------------------------|
| Total Kjeldahl nitrogen                              | U.S. EPA Method 351.1   |
| Ammoniacal nitrogen                                  | U.S. EPA Method 350     |
| Phosphate  | U.S. EPA Method 365     |
| Biological oxygen demand (5-day) (BOD <sub>5</sub> ) | U.S. EPA Method 405.1   |
| Chemical oxygen demand (COD)                         | U.S. EPA Method 410.1   |
| Total metals   | U.S. EPA Method 200.7   |
| 1 Otal Iliciais                                      | Cigi an 11 Montoe 2001. |

Additional Analyses. In addition to the aforementioned tests, the following tests were also conducted on groundwater and saturated zone soil samples:

| Microbial Enumerations        | Modified Spread Plate Method    |
|-------------------------------|---------------------------------|
| Microbial Stimulation Testing | IT Standard Operating Procedure |
| Minerals Analysis             | IT Standard Operating Procedure |
| Nutrient Compatibility        | IT Standard Operating Procedure |
| Nutrient Adsorption           | IT Standard Operating Procedure |
| Total Organic Carbon          | Modified U.S. EPA Method 415    |
| Oxygen Delivery Potential     | IT Standard Operating Procedure |

These tests provide additional information that is valuable in determining those physical, chemical, and microbiological properties that impact successful bioremediation. A brief description of each test is provided:

- Microbial enumerations are used to determine the population density of total bacteria and contaminant-degrading bacteria.
- Microbial stimulation testing is used to determine if the native bacterial population will respond favorably to oxygen and nutrient augmentation.
- Mineral analysis determines the calcium, magnesium, and iron content of site groundwater. These elements impact the efficiency of nutrient and oxygen transport.
- Nutrient compatibility examines the interaction of nutrients with groundwater and determines the proper nutrient addition rate.
- Nutrient adsorption evaluates the binding affinity of nutrients to aquifer materials.
   This relates to nutrient transport efficiency.
- TOC determinations provide insight on the ultimate mass of organic material that may
  be available for biodegradation in groundwater. This impacts the overall nutrient and
  oxygen requirements for the treatment system.
- Oxygen delivery potential determines the potential for efficient oxygen delivery through an aquifer using hydrogen peroxide.

#### 4.4.3.2 Microbial Inhibition Testing

The previously described microbial stimulation test will be used as a direct indicator of the inhibition of microbial activity by site samples. Microbial stimulation testing determines the response of native bacteria to nutrient and oxygen amendments. The response of microbes to this test is measured by increase in cell density. If the groundwater or soil contains toxic materials or is in some way inhibitory to microbial activity, the indigenous microbial population density fails to increase or is reduced. The microbial stimulation test should be independently conducted on groundwater and soil samples because many organic contaminants are found in higher concentrations in the soil. Analyzing only groundwater may give misleading results regarding the toxicity of the subsurface environment to microorganisms.

## 4.4.3.3 Batch Biodegradation Studies

Batch biodegradation studies were performed on site groundwater samples. The studies were conducted in computerized respirometers, which measure oxygen consumption by microorganisms and replace the oxygen to avoid oxygen limitations. This approach for monitoring microbial activity is not disruptive to the sample vessel and is not labor intensive.

Respirometric analysis of microbial activity provided a real-time indication of contaminant biodegradation. Using this approach, sampling points were chosen based on microbial activity rather than on an arbitrary time schedule. Also, because the vessels were sealed and remain sealed for most of the study, the loss of volatile organics was minimized.

A total of 15 treatment vessels were established. The treatment scheme is described in Table 1. Each treatment was conducted in triplicate. This experimental design demonstrated the biodegradation of target compounds in contrast to untreated and biologically inhibited controls. In addition to using indigenous microorganisms, the benefit of augmenting groundwater with activated municipal sludge and microorganisms isolated from site soil samples and grown on DEHP was evaluated. The use of alternative sources of microorganisms provided a means to increase the overall microbial density and diversity compared to groundwater alone. Activated sludge was collected from a local publicly owned treatment works (POTW) in Knoxville, Tennessee and a washing program was completed to deplete the residual dissolved carbon in the sludge. To ensure that each treatment contained the same initial concentration of target compounds, soil microorganisms were extracted from the soil and added to appropriate treatments in a minimal volume of contaminant-free buffer solution. Biological solids were added to a final solids density of 860 mg/L.

A single groundwater composite (18-L) was prepared from which all treatments were established. Triplicate analyses were performed on this composite. The composite groundwater sample was found to have a very low concentration of ethyl benzene, xylene, and DEHP; therefore, 3 mL of a

solution consisting of 1 part ethyl benzene, 1 part ortho-xylene, and 1 part DEHP was added to the 18-L composite sample and thoroughly mixed with the water to aid dissolution. An equal mass of water was removed from the bottom of the composite sample container and placed in each respirometer vessel. Nutrients and alternative microorganisms were added to appropriate treatments. Biologically-inhibited controls were treated with mercuric chloride. The untreated controls were placed in zero headspace vessels with no nutrient or microorganisms added and incubated in the dark at room temperature for the duration of the study. The remaining vessels were connected to the respirometer, maintained at constant temperature in the dark, and stirred continuously using a magnetic stir bar.

Oxygen consumption was continuously measured and recorded every 2 hours. Carbon dioxide production, pH, residual nutrients, and microbial density were determined at the end of the study. Contaminant and total organic carbon analyses were performed initially on the composite sample, on each vessel twice during the study as indicated by microbial respiration, and at the end of the study. The analytical scheme for the study is presented in Table 2. The study was terminated when microbial respiration became asymptotic.

#### 4.4.3.4 Soil Testing

Microbial Enumeration Testing. The density of heterotrophic and contaminant-degrading microorganisms was determined for soils used in the treatability studies. Although this phase of the study was in support of the biodegradation studies, it was also designed to support soil flushing results. The impact of biodegradation on the overall removal of contaminant mass was considered because this may significantly enhance the efficacy of soil remediation. Therefore, in support of the biodegradation studies, each soil sample was tested for pH, residual nutrients, and microbial stimulation as described in Section 4.2. Microbial population size and pH were also determined on representative aliquots of all four soils used in the soil flushing study prior to the initiation of flushing.

Initial Analytical Analyses. Samples received at the BAC from WESTON represent a range in contaminant levels for soils at the Carpenter site, as described in the sampling plan (Appendix A). To evaluate the efficacy of soil flushing as a potential remedial technology for site soils, soils with the highest concentration of contaminants were selected for use in the study. To determine the samples to be flushed, an initial analysis for DEHP from selected samples was conducted prior to initiating the study.

Aliquots of soil were selected from Shelby tubes retrieved from locations with reported high and medium concentrations of target compounds. Soil was equilibrated with an extraction solvent at a 1 to 5 (weight to volume) ratio for 12 to 18 hours. This slurry was extracted using sonication for 10 minutes prior to centrifugation. The supernatant was collected, concentrated, and analyzed for

DEHP using GC. Results from this initial screening were used to select the four samples used in the soil flushing treatability study.

Representative soil samples were collected from soil adjacent to each of the actual intact soil samples used in the soil flushing study. Soil could not be removed from the actual intact sample used in the flushing study. These analyses were, therefore, an initial estimate of the concentrations of VOCs and SVOCs for the intact samples used in the flushing study. Initial analyses included the following:

- VOCs
- SVOCs
- Hq •
- Water content.

Initial soil analysis data for mass balancing the soil flushing system could not be used. These soils were not chemically identical and therefore variability in the level of individual chemicals could be significant. Even soil within an intact sample will not be homogeneous with respect to the concentration of chemicals; therefore, all mass balancing efforts for VOCs and SVOCs within the soil flushing system were derived by back-calculating (the summing of chemical constituents in the leachates, final soil, and resin columns).

The flushing solution was potable water. The pH of the water was adjusted to match that of the site groundwater. The flushing solution was analyzed for the following:

- pH
- Specific conductance
- COD
- VOCs
- SVOCs.

## 4.4.3.5 Geotechnical Analyses

Geotechnical characterization of soils was determined on samples retrieved from the Carpenter site. Multiple depths for each of the three locations were sampled and, therefore, should provide some representative spatial distribution of the physical characteristics for the site.

To assess the range of characteristics that may be encountered at the site, selected representative samples were analyzed for the following geotechnical characteristics:

- Water content
- Particle size distribution
- Organic carbon content
- Atterberg limits
- Bulk density

- Specific gravity
- Hydraulic conductivity.

Samples were analyzed for the above-listed characteristics. Selection of soil samples for geotechnical testing depended on a number of factors including visual observation of each sample, the zones or horizons represented by each of the sampling depths, and the site locations for each set of samples or sample depths.

#### 4.4.3.6 Soil Flushing Studies

The procedure followed for the soil flushing study is a modified version of hydraulic conductivity test methods for soils: U.S. EPA 9100, Saturated Hydraulic Conductivity, Saturated Leachate Conductivity, and Intrinsic Permeability; American Society for Testing and Materials (ASTM) D 5084-90, Measurement of Hydraulic Conductivity of Saturated Porous Materials Using a Flexible Wall Permeameter; and U.S. Army Corps of Engineers EM 1110, Appendix VII: Permeability Tests. The procedure incorporates the use of a flexible wall permeameter (Figure 1) for conducting bench-scale soil flushing treatability studies on intact soil samples.

The flexible wall permeameter is a plexiglass cell designed to house an intact soil sample of a specific diameter and height. The sample, placed between a stainless steel pedestal and top cap, is encased in an inert material, confined in a latex membrane, and placed inside the cell. The cell is filled with water and a confining pressure applied. The sample is isolated from the cell water except for the confining pressure exerted on the sample. The pedestal and top cap are plumbed with Teflon® tubing to stainless steel inlet and outlet ports on the outside of the cell, allowing for flushing solution to enter and leachate to exit the sample.

## 4.5 Data Management

IT is committed to providing services for hazardous and toxic materials management that meet the technical and economic needs of our clients, satisfy regulatory requirements, and are commensurate with current state of the art. To provide direction for corporate operations so that they are performed in a controlled manner that can be demonstrated, IT has developed and implemented a formal Quality Assurance (QA) Program. This program, established in 1973, operates in compliance with the Code of Federal Regulations, 10 CFR 50, Appendix B; American National Standards Institute/American Society of Mechanical Engineers (ANSI/ASME) NQA-1 "Quality Assurance Program Requirements for Nuclear Facilities"; and current U.S. EPA guidelines and recommendations (e.g., QAMS-005/80, "Interim Guidelines and Specifications for Preparing Quality Assurance Project Plans"). This program establishes policies to facilitate the implementation of regulatory requirements and to provide internal means for control and review that ensure the work performed by IT complies with all requirements.

Site-specific QA/quality control (QC) procedures will be in accordance with the following documents:

- IT Engineering Operations QA Manual, Revision 1, July 1, 1987 and IT Analytical Services QA Manual, Revision 1, February 1, 1988.
- U.S. EPA, "Interim Guidelines and Specifications for Preparing Quality Assurance Project Plans," QAMA-005/80.
- IT Analytical Services QA Manual, Revision 1, February 1, 1988.

#### 4.5.1 Data Acquisition

Measuring and test equipment used in the field or laboratory was subject to a formal calibration program as described in Section 5.5, Equipment Calibration and Control, of the Environmental Projects Group Engineering Operations QA Manual. Calibration of laboratory equipment was documented in the project-specific laboratory notebook. Equipment that failed calibration was taken out of service and a Notice of Equipment Calibration Failure record was completed or the instrument was adjusted to permit correct operation. Records of equipment calibration are maintained in the project file located in Knoxville, Tennessee.

Sampling of all aspects of the treatability study was conducted by IT personnel. Sampling was performed in accordance with Section 5.0, Data Acquisition, of the Environmental Projects Group Engineering Operations QA Manual. Field samples were collected by WESTON personnel using IT Standard Operating Procedures for collection of microbiological samples and the IT Chain-of-Custody/Request for Analysis form.

All analyses were performed by either ITAS or the IT BAC laboratory in Knoxville, Tennessee. The methods used are listed in Table 3. QA/QC procedures prescribed in the U.S. EPA methods were used in all analyses. During the treatability study, total solids, total suspended solids, total dissolved solids, pH, ammoniacal nitrogen, ortho-phosphate, microbial enumerations, total organic carbon, and quantification of target compounds were performed by the IT BAC. Standard operating procedures and QA/QC requirements were followed. Records of all analyses were recorded in a bound laboratory notebook dedicated solely to this investigation.

## 4.5.2 Data Reduction, Validation, and Reporting

Data reduction, validation, and reporting were in accordance with the requirements contained in the following:

• IT Engineering Operations QA Manual, Revision 1, July 1, 1987, Modeling and Design.

 IT Analytical Services QA Manual, Revision 1, February 1, 1988, Section 10.0, Data Verification.

Numerical analysis was performed in accordance with the implementation methods described in the aforementioned documents, as amplified or modified by the following discussion of numerical/analysis activities of particular significance.

To provide evidence of satisfactory work performance and the basis for information transmitted to the U.S. EPA, numerical analysis and results were documented. Documentation may include calculations, computer programs, and associated input/output, logs, drawings, and tables. Analysis activities will be performed in a planned and controlled manner. Where appropriate, documented and approved work instructions were employed.

Calculations are legible and in a form suitable for reproduction, filing, and retrieval.

Documentation is such that a technically qualified individual can review and understand the calculations and verify the results. Calculations were performed on standard calculation paper or laboratory notebooks whenever possible. Data validation was performed in accordance with the IT OA Manual, Section 6.2.1, Calculations.

IT drawings were signed and dated by the draftsperson performing the work and the project manager who has checked the drawing. Revisions were noted on the drawing original with a revision number and brief note describing each revision. The note was signed and dated by the draftsperson performing the work and the project manager who has checked the revision.

Data are reported in tabular form as part of the overall project report. Validation of the report data complies with the IT Engineering Operations QA Manual, Section 9.0, Report Preparation.

#### 4.5.3 Internal QC Checks

QA audits and surveillances are performed in accordance with the requirements contained in the following:

- IT Engineering Operations Quality Assurance Manual, Revision 1, July 6, 1990. Section 11.0, Quality Assurance Audits.
- IT Analytical Services Quality Assurance Manual, Revision 1, February 1, 1988, Section 14.0, Quality Assurance/Quality Control Audits.

QA audits and surveillances are performed in accordance with the implementation methods described in the aforementioned documents, as amplified or modified by the following discussion of QA audits and surveillance activities of particular significance. Monthly surveillances of the IT

Knoxville central files are performed. Sampling and Analysis Plan (SAP) audits are performed quarterly.

#### 4.5.4 Corrective Action

Nonconformance identification, reporting, disposition, corrective action performance, verification, and acceptance will be in accordance with the requirements contained in the following:

- IT Engineering Operations Quality Assurance Manual, Revision 1, July 1, 1987, Section 8.0, Nonconformances/Corrective Action.
- IT Analytical Services Quality Assurance Manual, Revision 1, February 1, 1988, Section 13.0, Nonconformances and Corrective Action.

Nonconformance identification, reporting, disposition, corrective action performance, verification, and acceptance were performed in accordance with the aforementioned documents, as amplified or modified by the following discussion of nonconformances and corrective action activities of particular significance.

Items, services, or activities that do not meet SAP requirements or IT accepted standard practice are to be reported as nonconformances if they meet any of the following criteria:

- Are detrimental to quality if not corrected immediately and are not corrected immediately
- Are detrimental to quality if not ultimately corrected, are not corrected immediately, and are not tracked by a system that will ensure their timely correction
- Are of a nature such that others can benefit from learning of the circumstances so they
  can take appropriate preventive action for their activities, and a Nonconformance
  Report is judged the most appropriate way to communicate this information
- Pertain to items that should be segregated or visually identified to preclude inadvertent use
- Are significant conditions adverse to quality in which case they require reporting so
  that the cause of the condition can be determined and appropriate corrective action
  taken to preclude recurrence
- Are equipment that fails calibration or becomes impaired during use and is not immediately repaired.

Nonconformance Reports prevent the installation of items, the use of services, or the performance of activities that are detrimental to quality and provide for their correction. Nonconformance Reports also disseminate "lessons learned" information so appropriate preventive measures can be taken by others. And finally, Nonconformance Reports can provide invaluable information

identifying the need for, and triggering the initiation of, corrective actions to preclude recurrence of significant conditions adverse to quality.

Each nonconformance affecting quality is documented, normally by the person that identifies or creates it. For this purpose, a Nonconformance Report form or audit report is used as appropriate. The project manager determines the appropriate corrective measures to be taken and obtains the concurrence of the QA Officer (QAO).

Nonconformances identified within IT analytical laboratories are reported in accordance with Section 13.0, Nonconformances and Corrective Action, of IT Engineering Operations (ITEO) QA Manual, except laboratory nonconformances identified during ITEO-performed audits, which are reported in accordance with Section 14.0 Quality Assurance/Quality Control, of the aforementioned document. The QAO is provided with copies of ITEO Nonconformance Reports and audit reports that identify nonconformances adversely affecting the quality of the data.

Nonconformances identified during audits conducted by or under the direction of the QAO are reported in accordance with Section 11.0, Quality Assurance Audits, of the IT Engineering Operations QA Manual.

All other nonconformances are to be reported on a Nonconformance Report. The person(s) identifying or creating a nonconformance is responsible for reporting it, unless assurance is provided that it will be reported by someone else and the issuance of the report is verified by the person identifying or creating the nonconformance.

#### 4.5.5 Data Collection

Data collected during execution of the study was recorded in a bound controlled laboratory notebook. Data generated from integrators and computerized instruments were printed and the resulting data sheets are kept with the project file. All data were verified and checked by a BAC scientist. Proof of verification is the dated signature of the checker at the bottom of each notebook page.

## 5.0 Results and Discussion

The results of the treatability study have been segregated into several sections describing the various aspects of the study.

#### 5.1 Biological Site Characterization

Chemical, physical, and microbiological characteristics of site samples were investigated. The following subsections discuss the results of this characterization.

#### 5.1.1 Site Characteristics

The chemical constituents present in site groundwater were evaluated during the initial phases of this study. The concentration of VOCs observed in five site samples is shown in Table 3. Ethyl benzene was the most prevalent compound with lower concentrations of toluene and benzene found in two wells. Table 4 shows the concentration of metals in a groundwater composite sample. The metals content was generally low with the exception of common soil metals, iron, calcium, magnesium and aluminum. Table 5 provides data on other groundwater characteristics.

#### 5.1.2 Site Bioassessment

In addition to the general site characteristics, soil and groundwater samples were subjected to a battery of tests designed to evaluate the potential for implementing a successful in situ bioremediation treatment strategy. Soil and groundwater samples were analyzed for several microbiological, physical, and chemical parameters known to impact bioremediation. Sample pH, soil moisture content, residual nutrient content, and groundwater mineral content were determined. The compatibility of nutrients with groundwater and the adsorption of nutrients by soil-groundwater slurries were also determined. Microbiological characteristics such as the population density of aerobic heterotrophic bacteria, hydrocarbon-degrading bacteria, and DEHP-degrading bacteria were quantified in each sample. For the purpose of this evaluation, hydrocarbon degraders are those organisms capable of using benzene, ethyl benzene, toluene, and xylene as their only carbon and energy source. The response of bacteria to nutrient and oxygen stimulation was also evaluated. The ability to transport oxygen was evaluated by examining the stability of hydrogen peroxide in soil-groundwater slurries and the potential for sustaining a high dissolved oxygen content in the sample slurries.

## 5.1.2.1 pH

The pH of each site soil and groundwater sample is shown in Tables 6 and 7, respectively. The optimal pH for bioremediation is generally accepted to be within the range of 6 to 8. A pH outside of this range may reduce microbial metabolism and biodegradation. The soil pH ranged from 5.6 to 7.7. All of the soil samples except one had a pH within the preferred range. The groundwater

sample had pH 6.4. Adjustment of the pH will not be necessary initially. Periodic monitoring will determine if adjustments become necessary as bioremediation proceeds.

#### 5.1.2.2 Nutrient Analysis

The nitrogen as ammonia and ortho-phosphate content of site soil and groundwater samples is shown in Tables 6 and 7. The nitrogen content of soil ranged from nondetectable to moderate. Ammoniacal nitrogen in groundwater was low. The phosphate content in the soil was generally high but nondetectable in the groundwater. Therefore, nutrient additions will be necessary for successful bioremediation.

Although the phosphate content of the soil was high and may be adequate for bioremediation, the availability of phosphate for use by soil microbes is difficult to predict because of the strong binding affinity between soil and phosphate. Periodic monitoring will determine the utilization of nitrogen and phosphate and also aid in establishing the frequency of nutrient augmentation.

## 5.1.2.3 Total Organic Carbon Analysis of Groundwater

The total organic carbon content of the groundwater was 17 to 18 mg/L (Table 8). This value represents low organic carbon loading in the groundwater.

#### 5.1.2.4 Mineral Analyses

The calcium, magnesium, and iron content of groundwater was evaluated because these elements can catalyze chemical reactions that may interfere with aquifer bioremediation. The divalent cations, calcium and magnesium, react with phosphate to form an insoluble precipitate. In the case of severe phosphate precipitation, wells and geological formations can become clogged.

Hydrogen peroxide is frequently used as an oxygen source for aquifer bioremediation. As hydrogen peroxide decomposes, oxygen is liberated. Soluble iron can react with hydrogen peroxide to form insoluble iron oxides and hydroxides. Because the rate at which peroxide decomposes is related to its ultimate efficiency in transporting oxygen through an aquifer, iron can limit oxygen movement.

The calcium, magnesium and iron content of site groundwater is shown in Tables 4 and 7. The concentration of calcium and magnesium was moderate; the iron content was high. Nutrient precipitation due to dissolved minerals may be a problem due to the moderate calcium content of the groundwater. Hydrogen peroxide decomposition is difficult to predict based on iron measurements alone because soil composition and the organic content of the soil also plays a significant role in hydrogen peroxide decomposition.

### 5.1.2.5 Nutrient Compatibility

In order to determine the interaction between calcium, magnesium, and phosphate, a nutrient compatibility test was performed. The microbial nutrient blend Restore® 375 contains tripolyphosphates that act as calcium and magnesium chelators; therefore, by adding more Restore® 375, less precipitation occurs due to the increased concentration of the chelator.

The initial addition of Restore® 375 (10,000 mg/L) was based on the chelating power of the nutrient blend and the mineral content of the groundwater. The groundwater was observed for the formation of precipitate following the first addition of nutrients. If precipitation occurred, additional doses of nutrient were applied until the precipitate redissolves.

The groundwater composite received the first nutrient amendment of 10,000 mg/L with no precipitation or cloudiness. These results indicate that nutrient amendment to the groundwater can be accomplished with no anticipated complications.

#### 5.1.2.6 Nutrient Adsorption

Soils generally have an affinity for ammonium and phosphate. The binding capacity of a soil impacts the ability to transport nutrients away from the injection area. The data in Figure 3 show the amount of ammonia, phosphate, and chloride bound to a composite of the soil samples after each of three nutrient additions. The results indicated that after the initial addition of nutrients, approximately 40 percent of the ammonium, 80 percent of the added phosphate, and none of the chloride bound to the soil. The adsorption of ammonium and phosphate to the soil was initially high but showed of signs of saturation and reduced adsorption with each subsequent addition of nutrients.

The results of the nutrient adsorption test suggest that significant amounts of the added nutrients will initially bind to the soil. However, binding capacity of the soil will saturate as nutrients are added, thus permitting nutrient transport through the aquifer.

## 5.1.2.7 Hydrogen Peroxide Stability/Oxygenation Potential

Transport of oxygen through an aquifer is usually the limiting factor for in situ bioremediation. Oxygenation has been achieved successfully by applying hydrogen peroxide. When hydrogen peroxide decomposes, the resulting oxygen may be present as dissolved oxygen or as free oxygen gas. The movement of oxygen through an aquifer depends on the persistence of hydrogen peroxide and the level of dissolved oxygen in the water over time.

Figure 4 tracks the persistence and concentration of dissolved oxygen in a 9 parts water to 1 part soil slurry composed of site samples. The slurry was initially amended with about 868 mg/L hydrogen peroxide. Oxygen concentration was quantitatively tracked until changes in the oxygen

distribution between gaseous and dissolved phases became asymptotic. The data in Figure 4 show the amount of oxygen lost to the gas phase and the amount of dissolved oxygen defined as dissolved oxygen plus residual hydrogen peroxide.

Hydrogen peroxide at 600 mg/L will decompose into approximately 280 mg/L of oxygen. After about 2 hours, the dissolved oxygen content of the slurry was about 100 mg/L, and release of gaseous oxygen, as a result of hydrogen peroxide decomposition, was becoming asymptotic. Continued monitoring for a total of 4.5 hours indicated that about 70 mg/L of dissolved oxygen was still present.

A second application of 600 mg/L hydrogen peroxide was made. Decomposition progressed at about the same rate observed for the first application. Overall, the results indicated moderate hydrogen peroxide stability. Oxygenation of the aquifer using hydrogen peroxide should be acceptable.

#### 5.1.2.8 Microbial Enumerations

The spread plate technique was used to determine the density of the microbial population in each sample. Total heterotrophic microbes were grown on dilute nutrient agar. Hydrocarbon-degrading microorganisms were cultured on a mineral salts medium with the sole carbon and energy source being benzene, ethyl benzene, toluene, and xylene vapor. Phthalate degraders were cultured on mineral salts medium containing DEHP as the sole carbon and energy source. Microbial enumerations are shown in Table 9. Based on practical experience, the precision of the plate count method is approximately 0.5 order of magnitude.

The heterotrophic, hydrocarbon-degrading, and phthalate-degrading microbial population densities observed in soil samples were high. The microbial density in groundwater samples was moderate to high.

### 5.1.2.9 Microbial Stimulation

The response of native microorganisms to oxygenation and nutrient augmentation is an important factor in determining the likelihood of enhancing biodegradation. Table 10 shows the results of microbial stimulation tests. Because stimulation was measured as growth of bacteria, the plate count method was used to quantitate the level of microbial activity induced by oxygenation and nutrient augmentation. Increase in microbial population density greater than 0.5 order of magnitude is considered to be a positive response. As indicated in Table 10, heterotrophs, hydrocarbon-degrading bacteria, and phthalate-degrading bacteria in most samples responded to oxygenation and nutrient amendment by increasing in cell density by greater than 0.5 order of magnitude. Heterotrophs in sample ST-2A failed to respond to treatment, hydrocarbon degraders showed a weak response, but phthalate degraders responded very well to both treatments. Bacteria

in groundwater (Table 10) responded favorably to stimulation as indicated by growth in the presence of oxygen and nutrients.

In general, the results of the microbial stimulation tests indicated that under appropriate conditions, microbes present within the impacted soils and aquifers should respond favorably to nutrient and oxygen amendment. Nutrients enhanced the growth of bacteria in all cases except heterotrophs and hydrocarbon degraders in sample ST-2A.

#### 5.2 Biotreatability Study Results

The results of each evaluation made during the groundwater batch biotreatability study are described in the following sections.

## 5.2.1 Reduction in Target Compound Concentration During Treatment

The loss of ethyl benzene, xylene, and DEHP was determined at four points during the batch groundwater biotreatability test (Figure 5). The results indicate that biological treatment contributed to the removal of target compounds from biologically active treatments. The biologically inhibited treatment (GW, HgCl & Ntr) indicates the loss of target compounds through physical processes such as volatilization and adsorption to the reaction vessel. The vapor pressures of ethyl benzene and xylene are high enough that an appreciable amount of these compounds should volatilize and be lost when the treatment vessels were opened for sampling. The results shown in Figure 5 suggest this occurred. The vapor pressure of DEHP is so low that volatilization is not a major contributor to the abiotic loss of DEHP from the treatment vessels. The zero headspace treatment represents the absolute change attributable to experimental and analytical artifacts in a completely untreated control. A 45 percent loss of ethyl benzene, 33 percent loss of xylene, and 18 percent loss of DEHP was observed in this control treatment based on the analysis performed by the BAC laboratory. Ethyl benzene and xylene removal in all biologically active controls was greater than 99.9 percent. Volatilization could account for 80 percent of the ethyl benzene loss and 75 percent of the xylene loss. DEHP removal in the two treatments using indigenous microbes was approximately 99.7 percent. DEHP removal by the activated sludge treatment was only 86 percent.

The significance of target compound removal was evaluated using ANOVA and the F Test. Complementary statistical analyses were performed based on BAC analysis of each treatment and target compound analysis performed according to U.S. EPA protocols at an IT Analytical Services laboratory. The results of the statistical comparisons of each treatment are shown in Tables 11 and 12.

The results are the same using either analytical method for ethyl benzene and xylene. For these two compounds, a significant difference at the 95 percent confidence level was observed between

the biologically active treatments and the abiotic and the untreated conditions. There was also a significant difference between the abiotic treatment and the zerohead space control. This difference is attributed to volatilization in the abiotic treatment that was opened two more times than the zero headspace treatment. No difference in the removal of ethyl benzene and xylene was observed among any of the biologically active treatments. These results strongly suggest that the biological component of the treatment process contributed to the removal of ethyl benzene and xylene.

The same statistical process was used to evaluate the removal of DEHP from each treatment. The results vary slightly depending on the method of DEHP analysis. Biological treatment was shown to significantly contribute to DEHP removal at the 95 percent confidence level. Using BAC analytical data for DEHP removal, a significantly greater DEHP removal was detected in groundwater treated with indigenous microbes compared to the treatment augmented with activated sludge. No difference was detected between the native groundwater and groundwater augmented with DEHP degrading bacteria isolated from site groundwater. Using DEHP data generated by U.S. EPA methodology, no difference was detected among the biologically active treatments.

## 5.2.2 Biodegradation Rate Constants and Half-Lives

The biodegradation rate constant was determined from the target compound removal data. Figure 6 indicates the degradation rate constant and half-life for DEHP in each treatment. The biologically active treatments had larger rate constants and shorter half-lives than the mercury treated control. The treatments containing indigenous microbes had much shorter half-lives for DEHP than the activated sludge treatment.

The removal data for ethyl benzene and xylene were not appropriate for regression analysis of rate constants and half-lives; therefore, the initial concentration and the concentration at the first sampling point were used. In the case of ethyl benzene, the degradation rate exceeded the sampling frequency; therefore, the concentration was nondetectable at the first sampling point. The estimated biodegradation rate constant for ethyl benzene is  $\geq$  -0.2 hrs-1 with a half-life of  $\leq$  3.3 hrs based on complete removal within 43 hours. The biodegradation rate constant and half-life were also calculated for xylene using the initial concentration and the concentration at the first sampling point. The rate constant for the activated sludge treatment was -0.2 hrs-1 with a half-life of 3.3 hours. The rate constants for the two treatments containing indigenous microbes were essentially the same with a rate constant of -0.02 hrs-1 and a half-life of 37 to 39 hours.

## 5.2.3 Total Dissolved Organic Carbon Content During Treatment

The reduction in TOC during the test period is shown for each treatment in Figure 7. The general trends in TOC removal are consistent with the removal of target compounds in Figure 5. The TOC in the activated sludge augmented treatment is higher than for other biologically active treatments due to the TOC contributed by decaying sludge. The TOC in the DEHP-degrading bacteria

augmented treatment (GW, Soil, & Ntr) was slightly higher than the groundwater (GW & Ntr) treatment for the first two sampling points. At later points in the test, the TOC of these two treatments was essentially the same. The initially higher TOC in the GW, Soil & Ntr treatment was attributed to soluble organic carbon associated with the bacteria used to augment this treatment. Volatile suspended solids analysis of these treatments supports this interpretation. As indicated in Figure 8, the volatile solids content of these two treatments quickly equalized to the same value and remained very similar for the rest of the study. The microbial density of these treatments, reported in Figure 9, also supports population equilibration to the same level.

## 5.2.4 Respiration During Treatment

Figure 10 tracks microbial respiration in each treatment except the zero headspace control. The figure shows the mean  $\pm$  one standard deviation at each point collected during the study. Because the data collection frequency was 2 hrs, data were plotted as a line rather than as individual points. The results indicate that the activated sludge augmented treatment consumed over two times more oxygen than the other biologically active treatments; however, target compound loss was not enhanced as shown in Figure 5 and Tables 11 and 12. This treatment represents an unfavorable alternative because much more oxygen was required to achieve the same result. Oxygen consumption by the treatment augmented with DEHP degraders was slightly greater than the unaugmented groundwater treatment. The effectiveness of the mercury treatment for inhibiting biological activity was demonstrated by the virtual absence of oxygen consumption by these treatments.

## 5.2.5 Inorganic Carbon Production During Treatment

The production of inorganic carbon (CO<sub>2</sub>) complements the consumption of oxygen. However, absolute comparisons of oxygen consumed and carbon dioxide produced are difficult because some organic carbon is converted into biomass rather than being mineralized. Table 13 shows the amount of oxygen consumed, the total inorganic carbon (TIC) (carbon dioxide) generated, and the TOC consumed during the study. Observed oxygen consumption was slightly higher than expected based on TIC produced and TOC consumed for the groundwater treatment and the DEHP degrader augmented treatment; however, the slight difference in observed and expected oxygen consumption suggests that both treatments efficiently mineralized TOC. These results are in contrast to those observed for the activated sludge treatment which was much less efficient in mineralizing TOC.

## 5.2.6 Microbial Density During Treatment

Figure 9 tracks the microbial density of heterotrophs and phthalate-degrading bacteria at the start of treatment and at the end. The microbial density in the groundwater and nutrient treatment increased during treatment. The microbial density of the activated sludge and the phthalate-degrader amended treatments decreased during treatment. The abiotic control contained no bacteria and the

zero headspace treatment had a low stable heterotroph population. The phthalate degrader density decreased during incubation.

Total suspended solids (TSS) and volatile suspended solids (VSS) analyses were performed on each treatment. The average TSS and VSS for each treatment are shown in Figures 11 and 8, respectively. Solids analyses generally confirmed the microbial density determinations shown in Figure 9. However, the solids data for the abiotic control indicated an increase in TSS and VSS. This was due to a visibly noticeable precipitate caused by the presence 500 mg/L HgCl<sub>2</sub>. Precipitates were not observed in any other treatment.

Volatile solids for the groundwater and phthalate degrader-amended groundwater treatments had equalized by the end of the study. This observation suggests that the addition of extra bacteria will not have a long-term effect on the native population density. Similarly, the volatile solids content of the activated sludge treatment gradually fell after the first few hours of treatment.

### 5.2.7 Nutrient Utilization During Treatment

Ammonia and ortho-phosphate concentrations in each treatment were examined. No appreciable reduction in nutrient concentration occurred during treatment (Figures 12 and 13).

#### 5.2.8 Changes in pH During Treatment

Change in system pH was measured during treatment as indicated in Figure 14. Data represent the mean of the measurements taken from the triplicates of each treatment. The abiotic control and the zero headspace control had an essentially constant pH. The activated sludge treatment displayed an obvious acidification while the other biologically active treatments became slightly alkaline.

## 5.3 Soil Flushing Data Analysis and Interpretation

## 5.3.1 Sample Selection

Eighteen polyethylene sleeves from Shelby tube samples were received by the IT laboratory for use in a soil flushing treatability study. All samples were collected on February 7, 1992 by WESTON personnel and are referenced as to exact location and depth within the site (Appendix A). All sleeves were visually inspected so as to characterize the samples based on color, texture, and integrity (whether the sample was intact or disturbed). Based on this inspection and the sampling information, individual samples were selected for either geotechnical testing or soil flushing. Samples not considered to be suitable for either of these tests were not used in the study. All samples and their fate are listed in Table 14.

#### 5.3.2 Initial Phthalate Analysis

Soil samples to be used in the soil flushing treatability study were selected based on an initial analysis of the phthalate concentration in selected samples. A 50 g aliquot of soil was collected from the top part of each sleeve for analysis. Sample preparation and analysis was conducted according to standard operating procedures. Selection was based on sample integrity, sample depth, estimated phthalate concentration (determined during initial site characterization), and site location. Based on these criteria, five Shelby tube samples were selected from the 18 samples received by the laboratory.

The samples selected were ST-4D (12 to 36 inches), ST-3B (0 to 24 inches), ST-3A (24 to 48 inches), ST-2A (72 to 96 inches) and ST-2B (48 to 72 inches). The phthalate concentration for each sample was 0.04, 0.6, 11.1, 19.0 and 0.3 g/L, respectively. Based on this information, samples, ST-2A, ST-3A, and ST-4D were selected for use in the soil flushing treatability study. The ST-3A sample contained sufficient soil to be used in two cells of the flushing study.

#### 5.3.3 Geotechnical Characteristics

#### 5.3.3.1 Shelby Tube Samples

Six Shelby tube samples were selected from the remaining polyethylene sleeves for geotechnical characterization. Site areas that were targeted for geotechnical evaluation were those areas with low to medium levels of contamination. Samples selected represented four depth intervals, 0 to 24, 24 to 48, 48 to 72 and 72 to 96 inches. The vertical cross section of the soil profile depicted by these samples selected was considered representative of the geotechnical characteristics of the site. Geotechnical characteristics of the six selected soil cores are described in Table 15.

## 5.3.3.2 Soil Flushing Samples

Geotechnical characteristics of the soil samples used in the four flexible wall permeameter cells are given in Table 16. The three Shelby tube samples used in the soil flushing study are as follows:

| Cell 1 | ST-2A (72 to 96 inches)                  |
|--------|--|
| Cell 2 | ST-3A (24 to 48 inches) (top section)    |
| Cell 3 | ST-3A (24 to 48 inches) (bottom section) |
| Cell 4 | ST-4D (12 to 36 inches).                 |

An important geotechnical characteristic used in evaluating the flushing efficiency of a soil sample is the volume of voids. The total volume of voids in cubic centimeters (cc) for a given sample is described as one pore volume for that sample. This characteristic is used in describing the volume of flushing solution passing through the sample for a given flushing time period. The pore

volumes for the soils used in the flushing study ranged from a low of 53.9 cc for Cell 1 soil to a high of 153.7 cc for Cell 3 soil.

# 5.3.4 Initial Chemical Characterization of the Soil Flushing Samples Soil was collected from each of the Shelby tube sleeves selected for use in the flushing study. Samples were taken from a composite of the soil adjacent to the actual soil sample used in each of the cells. Because soil cannot be collected directly from the actual sample used in a cell, the homogenized soil sample provided an estimate of the analyte concentration for the soil in each cell.

These results are not intended to represent the actual starting concentration of analytes in the soil within the treatment cells.

All three soils were initially analyzed by GC/MS for priority pollutant volatile organic compounds VOCs (U.S. EPA SW-846 method 8240) and priority pollutant semivolatile organic compounds SVOCs (U.S. EPA SW-846 method 8270). A library search was conducted on all samples for other volatile and semivolatile nontarget analytes. TOC by U.S. EPA Methods 415.1 and 9060 was determined for each soil. Table 17 contains a listing of detectable analytes and their concentrations in each of the three soil samples selected for the flushing study.

The soil used in Cell 1 showed the highest concentration of VOCs and, in particular, the two targeted VOCs, ethyl benzene (130 mg/kg) and total xylenes (570 mg/kg). Total xylenes were detected Cells 2 and 3 and Cell 4 soils at 11 and 2.2 µg/kg, respectively. No ethyl benzene was detected in these two soils.

Phthalate concentrations in all three soils were very high, especially DEHP, the targeted SVOC in this study. The concentrations of DEHP in Cell 1, 2 and 3, and 4 soils were 1900, 650 and 840 mg/kg, respectively. The total phthalate concentrations equalled 0.2, 0.06, and 0.1 percent for Cells 1, 2 and 3, and 4, respectively. Few other priority pollutant SVOCs were detected in the initial soil analysis.

#### 5.3.5 Flushing Solutions and Flow Characterization

#### 5.3.5.1 Flushing Solutions

The soil flushing study was divided into two phases based on changes in the flushing solutions used for each phase. In Phase I potable water was used as the flushing solution in all four cells. The water was allowed to set open to the atmosphere for 24 hours prior to use in the flushing study. The pH of the water was 6.5.

In Phase II of the study, the flushing solution was changed. Initial results from the first leachate collection period indicated that little or no phthalates were eluted from the soil using just tap water.

Therefore, half way through the projected study period (30 days), the flushing solution changed for three of the four cells as described below:

| Cell 1 | 0.5 percent Brij 30/35 aqueous solution |  |
|--------|---|--|
| Cell 2 | $H_2O_2$ aqueous solution               |  |
| Cell 3 | Potable water                           |  |
| Cell 4 | 0.5 percent Tween 85 aqueous solution.  |  |

These flushing solutions were used throughout the 29 days of the second half of the soil flushing study.

#### 5.3.5.2 Flow Characterization

The flushing solution flow patterns for both study phases are illustrated in Figure 15. Flow volumes were determined gravimetrically and converted to volume using a 1.0 g/cc conversion factor. Both patterns show a linear flow volume with respect to time. A deviation in flow rate occurred after 20 days of operation. This slight depression in flow volume was a result of a temporary power failure to the peristaltic pump.

Tables 18 and 19 give individual and accumulative flow rates for all four samples during Phases I and II, respectively, of the study. Accumulative flow volumes for the 30 days of Phase I ranged from 26,926 mL for Cell 4 to 28,598 mL for Cell 1. During the 29 days of flushing in Phase II, accumulative flow volumes ranged from 26,230 mL for Cell 4 to 26,965 mL for Cell 3.

#### 5.3.6 Soil Flushing Results

#### 5.3.6.1 Phase I

The effectiveness of potable water as a flushing solution in removing VOCs and SVOCs from soils is illustrated in Tables 20 through 24. The results show the amount of each analyte detected in the leachate from each cell for two collection periods at Day 4 and Day 30. The amount of each analyte in a composite of all leachate from each cell is also presented, as well as the total amount each analyte collected during the entire Phase I flushing period. These values are derived from analyte concentration values (Appendix D) for each cell and collection period and represent the total amount of leachate collected for each flushing period (Table 18).

Water was effective in removing VOCs from each of the four soil samples. Relatively high amounts of ethyl benzene and total xylenes were removed during the first 4 days of soil flushing as compared to the amount removed during the final 3 days or the amount contained in the composite samples. It should be noted that the composite samples, although stored at 4°C following collection, were not analyzed until the conclusion of the study with storage times of 30 to 50 days.

This length of time may have resulted in the gradual loss of VOCs resulting in low to nondetectable levels in the composite samples.

Water was not effective in removing phthalates from soil. Little or no phthalates were detected in either of the collection period samples or the composite samples. The only phthalate that responded to flushing from all four samples was diethyl phthalate and this was only during the initial 4 day flushing period. No phthalates were detected in the leachates for the second collection period at 30 days, which was at the end of the Phase I flushing.

#### 5.3.6.2 Phase II

The results from the Day 4 leachate indicated that water was not an effective leaching agent for phthalates in soil. At this point, the water was amended with reagents to enhance the solubility of phthalates thus aiding in their removal from soil. Earlier efforts with two surfactants, Brij 30/35 and Tween 85, involving the removal of nonaqueous phase substances from soils, indicated that phthalates might also be effectively removed with a surfactant-amended flushing solution.

Therefore, these two surfactants were tested for the last 30 days in Phase II of the soil flushing study. Each surfactant was supplied via flushing solution to a single cell at a concentration of 0.5 percent weight to volume (Cell 1 - Brij 30/35, and Cell 2 - Tween 85). As a comparative measure, the flushing solution for Cell 3 was retained as potable water. In an effort to elucidate the effects of microbial enhancement on phthalate degradation, a hydrogen peroxide and nutrient-amended flushing solution was supplied to Cell 2 during Phase II of the study.

Brij 30/35 Flushing Solution. Table 20 shows the effect of a 0.5 percent Brij 30/35 surfactant amended flushing solution on the removal of VOCs and SVOCs from Cell 1 soil. Analysis of the leachate collected during the first 3 days of Phase II flushing indicated that ethyl benzene was the only VOC being flushed from the soil. A concentration of 1,800  $\mu$ g/L was detected in the 3-day leachate and 260  $\mu$ g/L was detected in the 29-day leachate. Once again the composite sample showed no detectable levels of any VOC.

The effect of a 0.5 percent Brij 30/35 flushing solution on removal of phthalates from soil is also illustrated in Table 20. The concentration of DEHP in the 3-day leachate was 61,000  $\mu$ g/L, and 3,200  $\mu$ g/L in the 29-day leachate. Analysis of the leachate composite collected during the interim flushing periods between days 3 and 26 of Phase II of the study indicated that a constant removal of this analyte was occurring. The concentration of DEHP in the composite leachate was 6,500  $\mu$ g/L.

The only other two phthalates that responded to a Brij 30/35 amended flushing solution were butylbenzylphthalate and di-n-octyl phthalate. However, they were detected at much lower concentrations than the DEHP.

Potable Water Flushing Solution. The continued use of potable water as a flushing solution during Phase II of the flushing study indicated a similar pattern as the Phase I results. Table 21 shows the effect of potable water in removing VOCs and phthalates from Cell 2 soil during the 29 days of flushing. Low concentrations of benzene and toluene were detected in the 3-day, 29-day, and composite leachates. However, potable water was ineffective in removing phthalates from soil. Only 2  $\mu$ g/L of diethyl phthalate was detected in the 29-day leachate.

H<sub>2</sub>O<sub>2</sub> and Nutrient Amended Flushing Solution. The effect of a H<sub>2</sub>O<sub>2</sub> and nutrient amended flushing solution on removing VOCs and SVOCs from Cell 3 soil is shown in Table 22. Very low concentrations of targeted VOCs were detected in the 29-day leachate, and only one phthalate, butylbenzylphthalate, was detected in the composite leachate. However, contaminant degradation by microbial avtivity was not expected to significantly increase the amount of analytes removed in the flushing solution. Approximately 30 mg/L of oxygen was consumed per pore volume of solution. This equates to a total consumption of 855 mg of oxygen over 30 days assuming a 3.8-hour retention time. Also, assuming that 5.3 mg of oxygen are required to mineralize 1 mg of organic carbon to carbon dioxide, 161 mg of organic carbon could have been mineralized during the 30-day treatment period. This is equivalent to 218 mg of DEHP. The absolute performance of the biologically enhanced cannot be determined because the true starting concentration of target compounds is not known.

Tween 85 Flushing Solution. Table 23 shows the effect of a 0.5 percent Tween 85 amended flushing solution on the removal of VOCs and SVOCs from Cell 4 soil. VOCs were not detected in the 3-day leachate and only low levels were detected in the 29-day and composite leachates. There were some responses from SVOCs to the Tween 85 amended flushing solution. Butylbenzylphthalate (190  $\mu$ g/L) and DEHP (490  $\mu$ g/L) were detected in the 3-day leachate during Phase II of the soil flushing study. However, only DEHP was detected in the 29-day leachate (360  $\mu$ g/L) and the composite leachate (38  $\mu$ g/L).

#### 5.3.7 Soil Flushing Effectiveness

The effectiveness of soil flushing should be evaluated based on a number of factors: (1) the initial and final concentrations of individual analytes in the soil; (2) the concentration of each analyte in the leachates as a function of time; and (3) the total amount of each analyte removed from the soil relative to the total number of pore volumes of flushing solution moving through the sample during the soil flushing study.

#### 5.3.7.1 Final Soil Analysis

The final soil analysis is given in Table 24. Comparison of these data with initial concentration data contained in Table 17 indicates that many of the analytes initially detected in the soil were not present in the soil at the conclusion of the flushing study.

VOCs were primarily detected in the Cell 1 sample (ST-2A) taken from the 72 to 96-inch soil depth. Of the eight VOCs (including two targeted VOCs, ethyl benzene and total xylenes) detected in the initial soil analysis, only ethyl benzene was detected in the final soil analysis.

There were 25 SVOCs detected in the initial soil analysis, seven of which were phthalic analytes. Only four SVOCs (three phthalic analytes) were detected in the final soil analysis. DEHP and dinoctyl phthalate were detected in all three soil samples at the conclusion of the soil flushing study.

It must be emphasized that the comparison between the initial and final soil concentrations is strictly a relative comparison, and in no way reflects a quantitative reduction in respective analytes. An estimate of the quantitative reduction in specific analytes can only be accomplished by mass balancing the analytes within the soil flushing system.

#### 5.3.7.2 System Mass Balancing

The total amount of each analyte removed during the soil flushing process is a product of the concentration of individual analytes in the leachate during each collection period and the total volume of leachate for that specific collection period. The total amount of VOCs and SVOCs removed from soils for each collection period during the two phases of the soil flushing study are given in Tables 25 through 28 for Cells 1 through 4, respectively.

The most interesting comparison within these tables is the difference in the amount (sum) of individual analytes collected during the Phase I and Phase II parts of the study. Most VOCs were removed from the soil during Phase I of the soil flushing process. However, data for Cell 1 (Table 25), which incorporated a 0.5 percent Brij 30/35 flushing solution, resulted in 50 percent more ethyl benzene being removed during Phase II of the study as compared to Phase I. It must be noted that this particular soil still had the highest concentration of ethyl benzene at the end of the study. Data in Table 28 also indicated that higher levels of benzene and toluene were removed from soil with the addition of a 0.5 percent Tween 85 flushing solution as compared to using just potable water.

The effect of adding surfactant amended flushing solutions on the total amount of phthalates removed from soil is clearly illustrated in Tables 25 and 28. Table 25 shows that butylbenzylphthalate, DEHP, and di-n-octyl phthalate were not removed with potable water; however, the addition of a 0.5 percent Brij 30/35 solutions removed 1,669, 369,928 and 1,221  $\mu$ g

of each analyte respectively. Even the use of a 0.5 percent Tween 85 flushing solution removed  $610\,\mu g$  of butylbenzylphthalate and 3,630  $\mu g$  of DEHP from the soil as compared to none being flushed using just potable water (Table 28).

Table 29 shows the total percent and amount of each analyte removed in the leachate during each phase of the soil flushing study. In addition, the total amount of VOCs adsorbed by the Tenax resin columns from the head space of the collection vessels and the amount of individual analytes still remaining in the soil at the conclusion of the flushing study are given. A mass balancing of individual analytes within the soil flushing system for this study is derived from summing these components to derive an original starting amount of each analyte within the soil at the beginning of the flushing study. A percent reduction in each analyte is then calculated based on this back-calculation method of mass balancing the system.

Almost 100 percent of all VOCs were removed from the soils in all four cells during the two phases of the soil flushing process. In many cases, over 50 percent of each analyte was removed during Phase I flushing using potable water. VOCs were extremely high in Cell 1 soil with over 17,000 µg (100 percent) removed during Phase I of the flushing study. In comparison, nearly 11,000 µg of ethyl benzene was flushed during the entire 60-day study, this equalled approximate 42 percent of the initial amount present in the soil.

Response of phthalates to soil flushing is also shown in Table 29. The most successfully flushing solution for removing phthalates, and specifically DEHP, was a 0.5 percent Brij 30/35 solution. Nearly 370 mg of DEHP (87 percent) was removed from Cell 1 soil during Phase II flushing using a surfactant-amended flushing solution as compared to none being flushed from the same soil use potable water during Phase I of the study. Sixty-four percent of butylbenzylphthalate (1,669  $\mu$ g) and 48 percent of di-n-octyl phthalate (1,221  $\mu$ g) were also flushed from the soil using the Brij 30/35 solution.

In contrast, only 2 percent of DEHP (3630  $\mu$ g) was flushed from Cell 4 soil using a 0.5 percent Tween 85 flushing solution. Phthalates were not detected in the leachate from Cell 2 and Cell 3 soils as a result of  $H_2O_2$ /nutrient amended and potable water flushing solutions. Much lower amounts of DEHP were detected in Cell 2 and Cell 3 final soils, which may have influenced its availability for flushing.

## Table 1 Batch Study Treatments L. E. Carpenter and Co., IT Project No. 408474

| Treatment                       | Nutrients and Oxygen | Number of Vessels |  |
|---------------------------------|----------------------|-------------------|--|
| Indigenous Microbes             | Yes                  | 3                 |  |
| Activated Sludge                | Yes                  | 3                 |  |
| Soil Microbes                   | Yes                  | 3                 |  |
| Biologically Inhibited          | Yes                  | 3                 |  |
| Zero Headspace Untreated Contro | l No                 | 3                 |  |

### Table 2 Batch Study Analytical Schedule L. E. Carpenter and Co., IT Project No. 408474

| Parameter                            |                     | Sampling Schedule   |
|--------------------------------------|---------------------|---|
| Total Solids (dissolved and suspend  | ded)                | $T_0, T_1, T_2, T_{final}$  |
| рН                                   |                     | T <sub>0</sub> , T <sub>1</sub> , T <sub>2</sub> , T <sub>final</sub> |
| Residual Nutrients (ammonia and ph   | nosphate)           | $T_0$ , $T_1$ , $T_2$ , $T_{final}$                                   |
| VOC and Base Neutrals                |                     | T <sub>0</sub> , T <sub>1</sub> , T <sub>2</sub> , T <sub>final</sub> |
| Total Organic Carbon                 | # 1                 | T <sub>0</sub> , T <sub>1</sub> , T <sub>2</sub> , T <sub>final</sub> |
| Total Inorganic Carbon               |                     | To and Tfinal   |
| Microbial Enumerations (total and ph | nthalate degraders) | T <sub>0</sub> and T <sub>final</sub>                                 |

Table 3
Volatile Organic Compounds in Groundwater
L. E. Carpenter and Co. IT Project No. 408474

| Compound                  |         |                 | Location        |         |                 |
|---------------------------|---------|-----------------|-----------------|---------|-----------------|
|                           | MW2     | MW3             | MW6             | MW14S   | MW15S           |
|                           | (µg/L)  | (μ <b>g/L</b> ) | (μ <b>g/L</b> ) | (µg/L)  | (μ <b>g/L</b> ) |
| Acrolein                  | ND (10) | ND (10)         | ND (10)         | ND (10) | ND (10)         |
| Acrylonitrile             | ND (10) | ND (10)         | ND (10)         | ND (10) | ND (10)         |
| Benzene                   | ND (5)  | 3               | 14              | ND (5)  | ND (5)          |
| Bromodichloromethane      | ND (5)  | ND (5)          | ND (5)          | ND (5)  | ND (5)          |
| Bromoform                 | ND (5)  | ND (5)          | ND (5)          | ND (5)  | ND (5)          |
| Bromomethane              | ND (10) | ND (10)         | ND (10)         | ND (10) | ND (10)         |
| Carbon Tetrachloride      | ND (5)  | ND (5)          | ND (5)          | ND (5)  | ND (5)          |
| Chlorobenzene             | ND (5)  | ND (5)          | ND (5)          | ND (5)  | ND (5)          |
| Chloroethane              | ND (10) | ND (10)         | ND (10)         | ND (10) | ND (10)         |
| 2-chloroethylvinyl ether  | ND (10) | ND (10)         | ND (10)         | ND (10) | ND (10)         |
| Chloroform                | ND (5)  | ND (5)          | ND (5)          | ND (5)  | ND (5)          |
| Chloromethane             | ND (10) | ND (10)         | ND (10)         | ND (10) | ND (10)         |
| Dibromochloromethane      | ND (5)  | ND (5)          | ND (5)          | ND (5)  | ND (5)          |
| 1,1-dichloroethane        | ND (5)  | ND (5)          | ND (5)          | ND (5)  | ND (5)          |
| 1,2-dichloroethane        | ND (5)  | ND (5)          | ND (5)          | ND (5)  | ND (5)          |
| 1,1-dichloroethene        | ND (5)  | ND (5)          | ND (5)          | ND (5)  | ND (5)          |
| Trans-1,2-dichloroethene  | ND (5)  | ND (5)          | ND (5)          | ND (5)  | ND (5)          |
| 1,2-dichloropropene       | ND (5)  | ND (5)          | ND (5)          | ND (5)  | ND (5)          |
| Cis-1,3-dichloropropene   | ND (5)  | ND (5)          | ND (5)          | ND (5)  | ND (5)          |
| Trans-1,3-dichloropropene | ND (5)  | ND (5)          | ND (5)          | ND (5)  | ND (5)          |
| Ethyl benzene             | 46      | 2000            | 13000           | ND (5)  | ND (5)          |
| Methylene chloride        | ND (5)  | ND (5)          | ND (5)          | ND (5)  | 1 (estimate     |
| 1,1,2,2-tetrachlorothane  | ND (5)  | ND (5)          | ND (5)          | ND (5)  | ND (5)          |
| Tetrachloroethene         | ND (5)  | ND (5)          | ND (5)          | ND (5)  | ND (5)          |
| Toluene                   | ND (5)  | 6               | 11              | ND (5)  | ND (5)          |
| 1,1,1-trichloroethane     | ND (5)  | ND (5)          | ND (5)          | ND (5)  | ND (5)          |
| 1,1,2-trichloroethane     | ND (5)  | ND (5)          | ND (5)          | ND (5)  | ND (5)          |
| Trichloroethene           | ND (5)  | ND (5)          | ND (5)          | ND (5)  | ND (5)          |
| Trichlorofluoromethane    | ND (5)  | ND (5)          | ND (5)          | ND (5)  | ND (5)          |
| Vinyl chloride            | ND (10) | ND (10)         | 3 (estimate)    | ND (10) | ND (10)         |

ND, nondetectable

## Table 4 Metals in Groundwater L. E. Carpenter and Co. IT Project No. 408474

| Metal     | Composite Water Sample |  |
|-----------|------------------------|--|
|           | (mg/L)                 |  |
| Aluminum  | 15.1                   |  |
| Antimony  | ND (0.03)              |  |
| Arsenic   | ND (0.05)              |  |
| Barium    | 2.8                    |  |
| Beryllium | ND (0.001)             |  |
| Cadmium   | 0.006                  |  |
| Calcium   | 65.6                   |  |
| Chromium  | 3.1                    |  |
| Cobalt    | 0,15                   |  |
| Copper    | 0.18                   |  |
| iron      | 427                    |  |
| Lead      | ND (0.03)              |  |
| Magnesium | 23.3                   |  |
| Manganese | 150                    |  |
| Nickel    | 2.4                    |  |
| Potassium | 5                      |  |
| Selenium  | 0.09                   |  |
| Silver    | ND (0.005)             |  |
| Sodium    | 12                     |  |
| Thallium  | ND (0.05)              |  |
| Vanadium  | 0.06                   |  |
| Zinc      | 0.23                   |  |

ND, nondetectable

### Table 5 Groundwater Parameters L. E. Carpenter and Co. IT Project No. 408474

| Parameter                       | Composite Sample |  |
|---------------------------------|------------------|--|
| <u></u>                         | (mg/L)           |  |
| Acidity as CaCO3                | 86               |  |
| Alkalinity as CaCO3             | 276              |  |
| Ammonia as N                    | 0.7              |  |
| Biochemical Oxygen Demand       | 14               |  |
| Chemical Oxygen Demand          | 57               |  |
| Total Kieldahl Nitrogen         | 3.3              |  |
| PH                              | 6.3              |  |
| Specific Conductance (µmhos/cm) | 490              |  |
| Total Dissolved Solids          | 290              |  |
| Total Solids                    | 500              |  |
| Total Suspended Solids          | 170              |  |

Table 6
Physical and Chemical Characteristics of Site Soil
L. E. Carpenter and Co., IT Project No. 408474

| рН  | Ammonium<br>(mg/kg)      | Phosphate<br>(mg/kg)                  | Soil Moisture<br>(percent)           |
|-----|--------------------------|---------------------------------------|--------------------------------------|
| 5.6 | <4                       | 80                                    | 8                                    |
| 7.7 | 63                       | 158                                   | 11                                   |
| 7.3 | 4.7                      | 221                                   | 13                                   |
| 7.4 | 19                       | 133                                   | 11                                   |
| 7.5 | 19                       | 139                                   | 8.                                   |
|     | 5.6<br>7.7<br>7.3<br>7.4 | (mg/kg)  5.6 <4 7.7 63 7.3 4.7 7.4 19 | (mg/kg)     (mg/kg)       5.6     <4 |

Table 7
Physical and Chemical Characteristics of Site Groundwater
L. E. Carpenter and Co., IT Project No. 408474

| Sample        | рH  | Ammonium<br>(mg/L) | Phosphate<br>(mg/L) | Calcium<br>(mg/L) | Magnesium<br>(mg/L) | Chloride<br>(mg/L) |
|---------------|-----|--------------------|---------------------|-------------------|---------------------|--------------------|
| Composite     | 6.4 | 4.7                | <0.5                | 61.6              | 33.3                | 45                 |
| Composite Dup | 6.4 | 4.4                | <0.5                | 61.6              | 29.1                | 45                 |

Table 8
Total Dissolved Organic Carbon in Site Groundwater
L. E. Carpenter and Co., IT Project No. 408474

| Sample       | Total Dissolved Carbon (mg/L) | Inorganic Carbon<br>(mg/L) | Organic Carbon<br>(mg/L) |
|--------------|-------------------------------|----------------------------|--------------------------|
| Composite    | 83                            | 66                         | 17                       |
| Composite Du | •                             | 65                         | 18                       |

Table 9
Enumeration of Heterotrophic and Hydrocarbon-Degrading
Bacteria
L. E. Carpenter and Co., IT Project No. 408474

| Sample        | Heterotrophs<br>(CFU/g) <sup>1</sup> | Hydrocarbon Degraders<br>(CFU/g) | Phthalate Degraders<br>(CFU/g) |
|---------------|--------------------------------------|----------------------------------|--------------------------------|
| Soil          |                                      |                                  |                                |
| ST-1B         | 4.3 X 10 <sup>6</sup>                | 1.6 X 10 <sup>6</sup>            | 1.5 X 10 <sup>6</sup>          |
| ST-2A         | 3.3 X 10 <sup>6</sup>                | 7.7 X 10 <sup>5</sup>            | 1.2 X 10 <sup>6</sup>          |
| ST-4A         | 4.7 X 10 <sup>6</sup>                | 1.5 X 10 <sup>6</sup>            | 1.1 X 10 <sup>6</sup>          |
| Composite     | 4.7 X 10 <sup>6</sup>                | 5.7 X 10 <sup>5</sup>            | 8.7 X 10 <sup>5</sup>          |
| Composite Dup | 4.3 X 10 <sup>6</sup>                | 1.9 X 10 <sup>6</sup>            | 8.2 X 10 <sup>5</sup>          |
| Groundwater   |                                      |                                  |                                |
| Composite     | 1.4 X 10 <sup>5</sup>                | 5.1 X 104                        | 8.2 X 10 <sup>3</sup>          |
| Composite Dup | 1.1 X 10 <sup>5</sup>                | 4.0 X 104                        | 6.2 X 10 <sup>3</sup>          |

<sup>1</sup>CFU/g, colony forming units per gram of dry soil or per mL for groundwater samples.

Table 10
Response of Indigenous Microorganisms to Oxygenation and Nutrient Amendment
L. E. Carpenter and Co., IT Project No. 408474
(Page 1 of 3)

| Sample        | Initial<br>(CFU/mL)   | Heterotrophs Oxygen (CFU/mL) | Oxygen and Nutrients<br>(CFU/mL) |
|---------------|-----------------------|------------------------------|----------------------------------|
| Soil          |                       | -                            |                                  |
| ST-1B         | 3.0 X 10 <sup>7</sup> | 5.2 X 10 <sup>7</sup>        | 1.4 X 10 <sup>8</sup>            |
| ST-2A         | 1.4 X 10 <sup>7</sup> | 2.0 X 10 <sup>7</sup>        | 1.1 X 10 <sup>7</sup>            |
| ST-4A         | 3.3 X 10 <sup>6</sup> | 3.0 X 10 <sup>7</sup>        | 9.4 X 10 <sup>7</sup>            |
| Composite     | 9.4 X 10 <sup>6</sup> | 1.7 X 10 <sup>8</sup>        | 4.0 X 108                        |
| Composite Dup | 7.4 X 10 <sup>6</sup> | 1.9 X 10 <sup>8</sup>        | 4.3 X 10 <sup>8</sup>            |
| Groundwater   |                       |                              |                                  |
| Composite     | 3.0 X 10 <sup>5</sup> | 8.8 X 10 <sup>6</sup>        | 3.5 X 10 <sup>7</sup>            |
| Composite Dup | 4.5 X 10 <sup>5</sup> | 2.3 X 10 <sup>7</sup>        | 8.9 X 10 <sup>7</sup>            |

**Table 10** (Page 2 of 3)

|               | Hydrocarbon-Degraders |                                   |                                  |  |
|---------------|-----------------------|-----------------------------------|----------------------------------|--|
| Sample        | Initial<br>(CFU/mL)   | Oxygen<br>(CFU/mL)                | Oxygen and Nutrients<br>(CFU/mL) |  |
| Soil          |                       |                                   |                                  |  |
| ST-1B         | 1.6 X 10 <sup>7</sup> | 6.9 X 10 <sup>7</sup>             | 3.5 X 10 <sup>8</sup>            |  |
| ST-2A         | 3.2 X 10 <sup>6</sup> | 1.1 X <del>1</del> 0 <sup>7</sup> | 7.5 X 10 <sup>6</sup>            |  |
| ST-4A         | 4.0 X 10 <sup>6</sup> | 1.3 X 10 <sup>7</sup>             | 3.1 X 10 <sup>8</sup>            |  |
| Composite     | 9.4 X 10 <sup>6</sup> | 1.7 X 10 <sup>8</sup>             | 4.0 X 10 <sup>8</sup>            |  |
| Composite Dup | 7.4 X 106             | 1.9 X 10 <sup>8</sup>             | 4.3 X 108                        |  |
| Groundwater   |                       |                                   |                                  |  |
| Composite     | 4.7 X 104             | 4.2 X 10 <sup>6</sup>             | 3.6 X 10 <sup>7</sup>            |  |
| Composite Dup | 6.0 X 104             | 4.4 X 106                         | 4.3 X 10 <sup>7</sup>            |  |

**Table 10** (Page 3 of 3)

|               |                       | Phthalate-Degrade     | re                            |
|---------------|-----------------------|-----------------------|-------------------------------|
| Sample        | Initial<br>(CFU/mL)   | Oxygen<br>(CFU/mL)    | Oxygen and Nutrients (CFU/mL) |
| Soil          |                       |                       |                               |
| ST-1B         | 1.4 X 10 <sup>7</sup> | 1.7 X 10 <sup>7</sup> | 1.3 X 10 <sup>7</sup>         |
| ST-2A         | 9.2 X 10 <sup>5</sup> | 4.8 X 10 <sup>7</sup> | 5.4 X 10 <sup>7</sup>         |
| ST-4A         | 1.5 X 106             | 6.6 X 10 <sup>6</sup> | 9.2 X 10 <sup>6</sup>         |
| Composite     | 1.4 X 10 <sup>6</sup> | 1.7 X 10 <sup>7</sup> | 2.0 X 10 <sup>7</sup>         |
| Composite Dup | 4.2 X 10 <sup>6</sup> | 6.6 X 107             | 4.1 X 10 <sup>7</sup>         |
| Groundwater   |                       |                       |                               |
| Composite     | 4.7 X 10 <sup>4</sup> | 1.1 X 10 <sup>6</sup> | 7.5 X 10 <sup>6</sup>         |
| Composite Dup | 3.7 X 104             | 1.1 X 10 <sup>6</sup> | 1.0 X 10 <sup>7</sup>         |

Table 11

ANOVA and Comparisons Among Means Based on BAC Lab Data

L.E. Carpenter and Co., IT Project No. 408474

| Ethylbenzene            | GW+Ntr     | GW+Sludge+Ntr | GW+Soil+Ntr | GW+HgCl+Ntr | Zero Headspace | ·····                |
|-------------------------|------------|---------------|-------------|-------------|----------------|----------------------|
|                         | 0          | 0             | 0           | 9.9         | 22.3           |                      |
|                         | 0          | 0             | 0           | 8.3         | 26.5           |                      |
|                         | 0          | 0             | 0           | 8.5         | 33.6           |                      |
| ΣΥ                      | 0.00       | 0.00          | 0.00        | 26.70       | 82.40          |                      |
| mean Y                  | 0.00       | 0.00          | 0.00        | 8.90        | 27.47          |                      |
|                         | 109.10     |               |             |             |                |                      |
|                         | 0.00       | 0.00          | 0.00        | 239.15      | 2328.50        |                      |
|                         | 2567.65    |               |             |             |                |                      |
|                         | 2500.88    |               |             |             |                |                      |
| GrandTotal^2/an         | 793.52     |               |             |             |                |                      |
| SStotal                 | 1774.13    |               |             | emer e      | ÷              |                      |
| SSgroup                 | 1707.36    |               |             |             |                |                      |
| SSwithin                | 66.77      |               |             |             |                |                      |
| ANOVA Table             | a= :       | 5.00          | n=          | 3.00        | F[0.05(4,10)]= | 3.48                 |
| Source of Variation     | d <i>f</i> | SS            | MS          | Fs          |                |                      |
| Among Groups            | 4.00       | 1707.36       | 426.84      | 63.93       | P≤0.05         |                      |
| Within Groups           | 10.00      | 66.77         | 6.68        |             | t.             |                      |
| Total                   | 14.00      | 1774.13       |             |             |                |                      |
|                         |            |               |             |             |                |                      |
| Comparisons, Planned    |            |               |             |             |                |                      |
| Source of variation     | SS         | df            | MS          |             |                | F(0.05(4.40)) 0.40   |
| Treatments              | 1707.36    | 4.00          |             |             | P≤0.05         | F[0.05(4,10)]= 3.48  |
| Killed vs Biological    | 237.63     | 1.00          |             |             | P≤0,05         | F[0.05(1,10)] = 4.96 |
| Untreated vs Biological | 2263.25    | 1,00          |             |             | P≤0.05         | F[0.05(1,10)] = 4.96 |
| Sludge vs Indigenous    | 0.00       | 1.00          |             |             |                | F[0.05(1,10)] = 4.96 |
| Sludge vs GW            | 0.00       | 1.00          |             |             |                | F[0.05(1,10)]= 4.96  |
| Sludge vs Soil          | 0.00       | 1.00          |             |             |                | F[0.05(1,10)] = 4.96 |
| GW vs Soil              | 0.00       | 1.00          |             |             |                | F[0.05(1,10)] = 4.96 |
| Untreated vs Killed     | 2025.62    | 1.00          |             |             | P≤0.05         | F[0.05(1,10)]= 4.96  |
| Within                  | 66.77      | 10.00         | 6.68        |             |                |                      |
| Total                   | 1774.13    | 14            |             |             |                |                      |

Table 11. Page 2 of 3

| o-Xylene                | GW+Ntr  | GW+Sludge+Ntr | GW+Soil+Ntr | GW+HgCl+Ntr | Zero Headspace |                      |
|-------------------------|---------|---------------|-------------|-------------|----------------|----------------------|
|                         | 0       | 0             | 0           | 13.1        | 25.7           |                      |
|                         | 0       | 0             | 0           | 11.3        | 30.6           |                      |
|                         | 0       | 0             | 0           | 11.4        | 37.6           |                      |
| Σy                      | 0.00    | 0.00          | 0.00        | 35.80       | 93.90          |                      |
| mean y                  | 0.00    | 0.00          | 0.00        | 1193        | 31.30          |                      |
| ΣΣΥ                     | 129.70  | •             |             |             |                |                      |
| Σ(y^2)                  | 0.00    | 0.00          | 0.00        | 429.26      | 3010.61        |                      |
| Σ(ΣΥ)^2                 | 3439.87 |               |             |             |                |                      |
| Σ(ΣY)^2/n               | 3366.28 |               |             |             |                |                      |
| GrandTotal^2/an         | 1121.47 |               |             |             |                |                      |
| SStotal                 | 2318.40 |               |             |             |                |                      |
| SSgroup                 | 2244.81 |               |             |             |                |                      |
| SSwithin                | 73.59   | ·             |             | <u> </u>    |                |                      |
| ANOVA Table             | a= :    | 5.00          | n=          | 3.00        | F[0.05(4,10)]= | 3.48                 |
| Source of Variation     | d∱      | SS            | MS          | Fs          |                |                      |
| Among Groups            | 4.00    | 2244.81       | 561.20      | 76.26       | P≤0.05         |                      |
| Within Groups           | 10.00   | 73.59         | 7.36        |             |                |                      |
| Total                   | 14,00   | 2318.40       |             | ·           |                |                      |
|                         |         |               |             |             | ,              |                      |
| Comparisons, Planned    |         |               |             |             |                |                      |
| Source of variation     | SS      | d <i>f</i>    | MS          | Fs          |                |                      |
| Treatments              | 2244.81 | 4.00          | 561.20      | 76.26       | P≤0.05         | F[0.05(4,10)]= 3.48  |
| Killed vs Biological    | 427.21  | 100           | 427.21      | 58.06       | P≤0.05         | F[0.05(1,10)]= 4.96  |
| Untreated vs Biological | 2939.07 | 1.00          | 2939.07     | 399.40      | P≤0.05         | F[0.05(1,10)]= 4.96  |
| Sludge vs Indigenous    | 0.00    | 1.00          | 0.00        | 0.00        | ns             | F[0.05(1,10)]= 4.96  |
| Sludge vs GW            | 0.00    | 1.00          | 0.00        | 0.00        | ns             | F[0.05(1,10)]= 4.96  |
| Sludge vs Soil          | 0.00    | 1.00          | 0.00        | 0.00        | ns             | F[0.05(1,10)] = 4.96 |
| GW vs Soil              | 0.00    | 1.00          | 0.00        | 0.00        | ns             | F[0.05(1,10)] = 4.96 |
| Untreated vs Killed     | 2511.86 | 1.00          | 2511.86     | 341.35      | P≤0.05         | F[0.05(1,10)] = 4.96 |
| Within                  | 73.59   | 10.00         | 7.36        |             |                |                      |
| Total                   | 2318.40 | 14            |             |             |                |                      |

Table 11. Page 3 of 3

| DEHP                    | GW+Ntr     | GW+Sludge+Ntr | GW+Soil+Ntr            | GW+HgCl+Ntr | Zero Headspace |                     |
|-------------------------|------------|---------------|------------------------|-------------|----------------|---------------------|
|                         | 0.1        | 2.5           | 0.07                   | 12.5        | 13.3           |                     |
|                         | 0.2        | 1.3           | 0.08                   | 8.58        | 14.7           |                     |
|                         | 0.2        | 3,5           | 0.05                   | 7.07        | 14             |                     |
| Σγ                      | 0.50       | 7.30          | 0.20                   | 28.15       | 42.00          |                     |
| mean y                  | 0.17       | 2.43          | 0.07                   | 9.38        | 14.00          |                     |
| ΣΣΥ                     | 78.15      |               |                        |             |                |                     |
| Σ(y^2)                  | 0.09       | 20.19         | 0.01                   | 279.85      | 588.98         |                     |
| Σ(ΣΥ)^2                 | 889.13     |               |                        |             |                |                     |
| Σ(ΣY)^2/n               | 870.00     |               |                        |             |                |                     |
| GrandTotal^2/an         | 407.16     |               |                        |             |                |                     |
| SStotal                 | 481.96     |               |                        |             |                |                     |
| SSgroup                 | 462.84     |               |                        |             |                |                     |
| SSwithin                | 19.12      |               |                        |             |                |                     |
| ANOVA Table             | a= 5       | .00           | Ú=                     | 3.00        | F[0.05(4,10)]= | 3.48                |
| Source of Variation     | d <i>f</i> | <b>SS</b> .   | MS                     | Fs          |                |                     |
| Among Groups            | 4.00       | 462.84        | 115.71                 | 60.50       | P≤0.05         | 1 - 1 - 1 - 1       |
| Within Groups           | 10.00      | 19.12         | 1.91                   |             |                |                     |
| Total                   | 14.00      | 481.96        | <u> </u>               |             |                |                     |
| Comparisons, Planned    |            |               |                        |             | •              |                     |
| Source of variation     | SS         | d <i>f</i>    | MS                     | Fs          | -              |                     |
| Treatments              | 462.84     | 4.00          | 115.71                 | 60.50       | P≤0.05         | F[0.05(4,10)]= 3.48 |
| Killed vs Biological    | 257.03     | 1.00          | 257.03                 | 134.40      | P≤0.05         | F[0.05(1,10)]= 4.96 |
| Untreated vs Biological | 580.89     | 1.00          | 580.89                 | 303.74      | P≤0.05         | F[0.05(1,10)]= 4.96 |
| Sludge vs Indigenous    | 17.69      | 1.00          | 1 <b>7</b> .6 <b>9</b> | 9.25        | P≤0.05         | F[0.05(1,10)]= 4.96 |
| Sludge vs GW            | 17.75      | 1.00          | 17.75                  | 9.28        | P≤0.05         | F[0.05(1,10)]= 4.96 |
| Sludge vs Soil          | 17.76      | 1.00          | 17.76                  | 9.29        | P≤0.05         | F[0.05(1,10)]= 4.96 |
| GW vs Soil              | 0.07       | 1,.00         | 0.07                   | 0.04        | ns             | F[0.05(1,10)]= 4.96 |
| Untreated vs Killed     | 323.86     | 1.00          | 323.86                 | 169.34      | P≤0.05         | F[0.05(1,10)]= 4.96 |
| Within                  | 19.12      | 10.00         | 1.91                   |             |                |                     |
| Total                   | 481.96     | 14            |                        |             |                |                     |

Table 12

ANOVA and Comparisons Among Means Based on EPA CLP Data

L.E. Carpenter and Co., IT Project No. 408474

| Ethylbenzene            | GW+Nuts    | GW+Sludge+Nuts GV | V+Soil+Nuts GW | +HgCl+Nuts | Zero Headspace |                      |
|-------------------------|------------|-------------------|----------------|------------|----------------|----------------------|
|                         | 0          | 0                 | 0              | 6.5        | 14             |                      |
|                         | 0          | 0                 | 0              | 3.6        | 15             |                      |
|                         | 0          | 0                 | 0              | 7.5        | 15             |                      |
| Σγ                      | 0.00       | 0.00              | 0.00           | 17.60      | 44.00          |                      |
| mean y                  | 0.00       | 0.00              | 0.00           | 5.87       | 14.67          |                      |
| ΣΣΥ                     | 61.60      |                   |                |            |                |                      |
| <br>Σ(y^2)              | 0.00       | 0.00              | 0.00           | 111.46     | 646.00         |                      |
| Σ(ΣΥ)^2                 | 757.46     |                   |                |            |                |                      |
| Σ(ΣΥ)^2/n               | 748.59     |                   |                |            |                |                      |
| GrandTotal^2/an         | 252.97     |                   |                |            |                |                      |
| SStotal                 | 504.49     | * 486 * 46.       |                | ***        | •              | -=                   |
| SSgroup                 | 495.62     |                   |                |            |                |                      |
| SSwithin                | 8.87       | -                 |                |            | .E             | -                    |
| ANOVA Table             | <b>a</b> = | 5.00              | n= 3.00        | 0          | F[0.05(4,10)]= | 3.48                 |
| Source of Variation     | d <i>f</i> | SS                | MS             | Fs         |                |                      |
| Among Groups            | 4.00       | 495.62            | 123.90         | 139.64     | P≤0.05         |                      |
| Within Groups           | 10.00      | 8.87              | 0.89           |            |                |                      |
| Total                   | 14.00      | 504.49            |                |            |                |                      |
|                         |            |                   |                |            |                |                      |
| Comparisons, Planned    |            |                   |                |            |                |                      |
| Source of variation     | SS         |                   | MS             | Fs         |                |                      |
| Treatments              | 495.62     | 4.00              | 123.90         | 139.64     | P≤0.05         | F[0.05(4,1.0)] = 3.4 |
| Killed vs Biological    | 103.25     | 1.00              | 103.25         |            | P≤0.05         | F[0.05(1,10)]= 4.9   |
| Untreated vs Biological | 645.33     | 1.00              | 645.33         |            | P≤0.05         | F[0.05(1,10)] = 4.9  |
| Sludge vs Indigenous    | 0.00       | 1.00              | 0.00           | 0.00       | ns             | F[0.05(1,10)]= 4.9   |
| Sludge vs GW            | 0.00       | 1.00              | 0.00           | 0.00       | ns             | F[0.05(1,10)]= 4.9   |
| Sludge vs Soil          | 0.00       | 1.00              | 0.00           | 0.00       | ns             | F[0.05(1,10)] = 4.9  |
| GW vs Soil              | 0.00       | 1.00              | 0.00           | 0.00       | ns             | F[0.05(1,10)] = 4.9  |
| Untreated vs Killed     | 542.08     | 1.00              | 542.08         | 610.91     | P≤0.05         | F[0.05(1,10)] = 4.9  |
| Within                  | 8.87       | 10.00             | 0.89           |            |                | •                    |
| Total                   | 504.49     | 14                |                |            |                |                      |

### Table 12. Page 2 of 3

| Xylene                  | GW+Nuts    | GW+Sludge+Nuts                          | GW+Soil+Nuts | GW+HgCI+Nuts | Zero Headspace |                      |
|-------------------------|------------|---|--------------|--------------|----------------|----------------------|
|                         | 0          | 0                                       | 0            | 5.1          | 15.8           |                      |
|                         | 0          | 0                                       | 0            | 3.55         | 15.9           |                      |
|                         | 0          | 0                                       | 0            | 9.39         | 26.9           |                      |
| Σγ                      | 0.00       | 0.00                                    | 0.00         | 18.04        | 58.60          |                      |
| mean y                  | 0.00       | 0.00                                    | 0.00         | 6.01         | 19.53          |                      |
| ΣΣΥ                     | 76.64      |   |              |              |                |                      |
| Σ(y^2)                  | 0.00       | 0.00                                    | 0.00         | 126.78       | 1226.06        |                      |
| Σ(ΣΥ)^2                 | 1352.84    | •                                       |              |              |                |                      |
| Σ(Σ <b>Y)^2/</b> n      | 1253,13    | }                                       |              |              |                |                      |
| GrandTotal^2/an         | 391.58     | }                                       |              |              |                |                      |
| SStotal                 | 961.27     | •                                       |              |              |                |                      |
| SSgroup                 | 861.55     | •                                       |              |              |                |                      |
| SSwithin                | 99.71      | - · · · · · · · · · · · · · · · · · · · |              |              |                |                      |
| ANOVA Table             | a=         | 5.00                                    | n=           | 3.00         | F[0.05(4,10)]= | 3.48                 |
| Source of Variation     | d <i>f</i> | SS                                      | MS MS        | Fs           |                |                      |
| Among Groups            | 4.00       | 861.55                                  | 215.39       | 21:60        | P≤0.05         |                      |
| Within Groups           | 10.00      | 99.71                                   | 9.97         | •            |                |                      |
| Total                   | 14.00      | 961.27                                  | <u> </u>     |              |                |                      |
|                         |            |   |              |              | ,              |                      |
| Comparisons, Planned    |            |   |              |              |                |                      |
| Source of variation     | SS         | d dj                                    | MS MS        | Fs           |                |                      |
| Treatments              | 861.55     | 4.00                                    | 215.39       | 21.60        | P≤0.05         | F[0.05(4,10)] = 3.48 |
| Killed vs Biological    | 108.48     | 3 1.00                                  | 108.48       | 10.88        | P≤0.05         | F[0.05(1,10)]= 4.96  |
| Untreated vs Biological | 1144.65    | 1.00                                    | 1144.65      | 114.80       | P≤0.05         | F[0.05(1,10)]= 4.96  |
| Sludge vs Indigenous    | . 0.00     | 1.00                                    | 0.00         | 0.00         | ns             | F[0.05(1,10)]= 4.96  |
| Sludge vs GW            | 0.00       | 1.00                                    | 0.00         | 0.00         | ns             | F[0.05(1,10)]= 4.96  |
| Sludge vs Soil          | 0.00       | 1.00                                    | 0.00         | 0.00         | ns             | F[0.05(1,10)]= 4.96  |
| GW vs Soil              | 0.00       | 1.00                                    | 0.00         | 0.00         | ns             | F[0.05(1,10)]= 4.96  |
| Untreated vs Killed     | 1036.17    | 7 1.00                                  | 1036.17      | 103.92       | P≤0.05         | F[0.05(1,10)]= 4.96  |
| Within                  | 99.71      | 10,00                                   | 9.97         | •            |                |                      |
| Total                   | 961.27     | 7 14                                    | 4            |              |                |                      |

Table 12. Page 3 of 3

| DEHP                    | GW+Nuts    | GW+Sludge+Nuts | GW+Soil+Nuts | GW+HgCl+Nuts | Zero Headspace |                                       |
|-------------------------|------------|----------------|--------------|--------------|----------------|---------------------------------------|
|                         | 0.11       | 1.8            | 0.15         | 11           | 4.2            |                                       |
|                         | 0.2        | 1              | 0.86         | 3.4          | 5              |                                       |
|                         | 0.18       | 1.9            | 0.27         | 3.4          | 9.2            |                                       |
| Σγ                      | 0.49       | 4.70           | 1.28         | 17.80        | 18.40          |                                       |
| mean y                  | 0.16       | 1.57           | 0.43         | 5.93         | 6.13           |                                       |
| ΣΣΥ                     | 42.67      |                |              |              |                |                                       |
| Σ(y^2)                  | 0.08       | 7.85           | 0.84         | 144.12       | 127,28         |                                       |
| Σ(ΣΥ)^2                 | 280.17     |                |              |              |                |                                       |
| Σ(ΣΥ)^2/n               | 226.46     |                |              |              |                |                                       |
| GrandTotal^2/an         | 121.38     | /              |              |              |                |                                       |
| SStotal                 | 158.79     |                |              |              |                |                                       |
| SSgroup                 | 105.07     |                |              |              |                |                                       |
| SSwithin                | 53.71      |                |              |              |                |                                       |
| ANOVA Table             | a= 5       | 5.00           | n=           | 3.00         | F[0.05(4,10)]= | 3.48                                  |
| Source of Variation     | d <i>f</i> | SS             | MS           | . Fs         |                |                                       |
| Among Groups            | 4.00       | 105.07         | 26.27        | 4.89         | P≤0.05         |                                       |
| Within Groups           | 10.00      | 53.71          | 5.37         |              |                |                                       |
| Total                   | 14.00      | 158.79         |              |              |                | · · · · · · · · · · · · · · · · · · · |
|                         |            |                |              |              | •              |                                       |
| Comparisons, Planned    |            |                |              |              |                |                                       |
| Source of variation     | SS         | df             | MS           | Fs           |                |                                       |
| Treatments              | 105.07     | 4.00           | 26.27        | 4.89         | P≤0.05         | F[0.05(4,10)]= 3.48                   |
| Killed vs Biological    | 100.96     | 1.00           | 100,96       | 18.80        | P≤0.05         | F[0.05(1,10)]= 4.96                   |
| Untreated vs Biological | 108.20     | 1.00           | 108.20       | 20.14        | P≤0.05         | F[0.05(1,10)] = 4.96                  |
| Sludge vs Indigenous    | 7.11       | 1.00           | 7.11         | 1.32         | ns             | F[0.05(1,10)]= 4.96                   |
| Sludge vs GW            | 7.35       | 1.00           | 7.35         | 1.37         | ns             | F[0.05(1,10)] = 4.96                  |
| Sludge vs Soll          | 7.30       | 1.00           | 7.30         | 1.36         | ns             | F[0.05(1,10)]=: 4.96                  |
| GW vs Soil              | -0.47      | 1.00           | -0.47        | -0.09        | ns             | F[0.05(1,10)]= 4.96                   |
| Untreated vs Killed     | 7.24       | 1.00           | 7.24         | 1.35         | ns             | F[0.05(1,10)]= 4.96                   |
| Within                  | 53.71      | 10.00          | 5.37         | •            |                |                                       |
| Total                   | 158.79     | 14             |              |              |                | <u> </u>                              |

Table 13
Respiration and Mineralization of Organic Carbon
L. E. Carpenter, IT Project No. 408474

| Parameter                              |              |                      | Treatment          |                    |                     |
|--|--------------|----------------------|--------------------|--------------------|---------------------|
|  | GW & Ntr (1) | GW, Sludge & Ntr (2) | GW, Soil & Ntr (3) | GW, HgCl & Ntr (4) | Zero Head Space (5) |
| Total Oxygen Consumed (mg/L)           | 397          | 1344                 | 474                | 13                 | ND                  |
| TIC Evolved (mg C//L)                  | 117          | 220                  | 140                | 64                 | ND                  |
| TIC Initial (mg C/L)                   | 56           | 57                   | <b>58</b>          | 57                 | .57                 |
| TIC Generated (mg C/L)(Evolved-Initial | 61           | 164                  | 82                 | .7                 | ND                  |
| TOC Consumed (mg C/L)                  | 55           | <b>36</b> °          | <b>62</b>          | NA                 | 0                   |
| Expected O2 based on TIC (mg/L)        | 325          | 874                  | 437                | 25                 | ND                  |
| Expected O2 based on TOC (mg/L)        | 293          | 194                  | 329                | NA .               | ND .                |

NA, not applicable

ND, not determined

Table 14
Cross Reference Chart for Soil Samples and Tests

| Split Spoon Sample | Shelby Tube          | Type of<br>Testing | Initial Phthalate<br>Analysis |
|--------------------|----------------------|--------------------|-------------------------------|
| Number             | Depth                |                    |                               |
| ST-1A              | 12 to 36             | Not Used           | No                            |
| ST-1A              | 36 to 60             | Not Used           | No                            |
| ST-1B              | 12 to 36             | Not Used           | No .                          |
| ST-1B              | 36 to 60             | Not Used           | No                            |
| ST-2A              | 24 to 48             | Geotechnical       | No                            |
| ST-2A              | 48 to 72             | Geotechnical       | No                            |
| ST-2A              | 72 to 96             | Soil Flushing      | Yes                           |
| ST-2B              | 24 to 48             | Geotechnical       | No                            |
| ST-2B              | 48 to 72             | Not Used           | Yes                           |
| ST-2B              | 72 to 96             | Geotechnical       | No                            |
| ST-3A              | 0 to 24              | Geotechnical       | No                            |
| ST-3A              | 24 to 48             | Soil Flushing      | Yes                           |
| ST-3B              | 0 to 24              | Not Used           | Yes                           |
| ST-3B              | 24 to 48             | Geotechnical       | No                            |
| ST-4A              | 12 to 36             | Not Used           | No                            |
|                    | 12 to 36             | Not Used           | No                            |
| ST-4B              |                      | Not Used           | No                            |
| ST-4C<br>ST-4D     | 12 to 36<br>12 to 36 | Soil Flushing      | Yes                           |

Table 15
Geotechnical Characteristics of Selected Soil Cores

| Sample Number                   | ST-2A   | ST-2A   | ST-2B   | ST-2B   | ST-3A   | ST-3B   |
|---------------------------------|---------|---------|---------|---------|---------|---------|
| Sample Depth (inches)           | 24-48   | 48-72   | 24-48   | 72-96   | 0-24    | 24-48   |
| Water Content (%)               | 6.5     | 15.1    | 6.6     | 5.6     | 10.2    | 40.2    |
| Average Specific Gravity (g/cc) | 2.82    | 2.56    | 2.95    | 2.83    | 2.63    | -2.51   |
| Organic Content (%)             | 5.3     | 8.0     | 4.2     | 3.8     | 3.4     | 4.3     |
| Ash Content (%)                 | 94.7    | 92.0    | 95.8    | 96.2    | 96.6    | 95.7    |
| Porosity                        | 34.6    | 67.1    | 49.3    | 43.3    | 38.5    | 46.8    |
| Liquid Limit                    | 26.0    | 30.0    | 24.0    | 23.0    | 20.0    | 30.0    |
| Plasticity Index                | 2.0     | NP      | NP      | 1.0     | 3.0     | NP      |
| Permeability (cm/s)             | 5.1 E-5 | 1.8 E-5 | 1.1 E-4 | 1.4 E-4 | 3.3 E-5 | 5.6 E-6 |
| Unit Weight Density (g/cc)      | 1.97    | 0.97    | 1.60    | 1.70    | 1.78    | 1.88    |
| Dry Density (g/cc)              | 1.85    | 0.84    | 1.50    | 1.61    | 1.62    | 1.34    |
| Total Organic Carbon (mg/kg)    | 18000   | 2300    | 3300    | 3500    | 4100    | 7000    |

NP denotes nonplastic

Table 16
Geotechnical Characteristics of Soil Samples Used in the Study

| Cell<br>Number | Soil<br>Weight<br>g | Soil<br>Volume<br>cc | Bulk<br>Density<br>g/cc | Volume of<br>Solids<br>cc | Volume of<br>Voids<br>cc | Void<br>Ratio | Pore<br>Volume<br>% |
|----------------|---------------------|----------------------|-------------------------|---------------------------|--------------------------|---------------|---------------------|
| Cell 1         | 520.13              | 246.5                | 2.11                    | 192.6                     | 53.9                     | 0.28          | 21.8                |
| Cell 2         | 358.81              | 280.8                | 1.28                    | 132.9                     | 147.9                    | 1.11          | 52.7                |
| Cell 3         | 491.50              | 335.7                | 1.46                    | 182.0                     | 153.7                    | 0.84          | 45.8                |
| Cell 4         | 535.45              | 323.3                | 1.66                    | 198.3                     | 125.0                    | 0.63          | 38.7                |

All calculations based on oven dry weight basis

Specific gravity estimated at 2.7 g/cc

Table 17
Concentration of Analytes in Initial Soil Samples Used in the Soil Flushing Study

| Cell Number   | Cell 1   | Cell 2/3 | Cell 4   |
|---|----------|----------|----------|
| Sample Identification                               | ST-2A    | ST-3A    | ST-4D    |
| Sample Depth (inches)                               | 72 to 96 | 24 to 48 | 12 to 36 |
| Chemical Analytes                                   | ug/kg    | ug/kg    | ug/kg    |
| Volatile Organic Compounds                          |          |          |          |
| Ethyl benzene                                       | 130000   | •        | 1        |
| 4-methyloctane                                      | 11000    | •        | ı        |
| Total xylenes                                       | 570000   | =        | 2.2      |
| 2,5-dimethylnonane                                  | 26000    | •        |          |
| Propylcyclohexane                                   | 40000    | ٠        | 1        |
| Ethylmethylbenzene                                  | 22000    | •        | ı        |
| 2-methylnonane                                      | 35000    | •        | 1        |
| Total cyclohexanes                                  | •        | 246.8    | •        |
| Semivolatile Organic Compounds                      |          |          |          |
| Butylbenzylphthalate                                | 1600     | 380      | 1500     |
| Bis(2-ethylhexyl)phthalate                          | 1900000  | 650000   | 840000   |
| Di-n-butylphthalate                                 | 11000    | 250      | 31000    |
| Diethyl phthalate                                   | 1600     | •        | 3700     |
| Dimethyl phthalate                                  | 380      |          | 760      |
| Di-n-octyl phthalate                                | 4700     | 1400     | 3000     |
| Total phthalates (additional)                       | 69000    | 16800    | 137000   |
| Anthracene  | •        | •        | 340      |
| Benzo(a)anthracene                                  | ı        | •        | 850      |
| Benzo(b)fluoranthene                                | 1        | •        | 640      |
| Benzo(k)fluoranthene                                | ľ        |          | 770      |
| Benzo(a)pyrene                                      | •        | •        | 720      |
| Chrysene  | •        | ı        | 890      |
| Fluoranthene  | •        | •        | 1500     |
| Indeno(1,2,3-cd)pyrene                              | •        | •        | 380      |
| Naphthalene   | •        | •        | 230      |
| Phenanthrene  | •        | •        | 1000     |
| Pyrene  | •        | •        | 1200     |
| Dibenz(a,h)anthracene                               | i        | •        | 250      |
| 4-Hydroxy-4-methyl-2-pentanone                      | 14000    | 14000    | 14000    |
| 2-Methyl-1-(1,1-dimethylethyl)-2-met propanoic acid | 16000    | 22000    | 5700     |
| Hexadecanoic acid                                   | 6700     | 1000     | 5700     |
| 2-Ethylhexyl diphenyl ester phosphoric acid         | 55000    | 8400     | 56000    |
| 2-Methylpropyl ester octadecanoic acid              | 9600     | •        | 9000     |
| Total alkanes                                       | 274400   | 1400     | 233800   |

Table 18
Leachate Volumes Collected During the Soil Flushing Study for Phase I

| Sample Period | 1      | 2      | 3     | 4     | .5     | 6      | 7      | 8      | 9      | 10     |
|---------------|--------|--------|-------|-------|--------|--------|--------|--------|--------|--------|
| Date          | 26-Mar | 30-Mar | 3-Apr | 6-Apr | 10-Apr | 11-Apr | 13-Apr | 15-Apr | 18-Apr | 21-Apr |
| Days          | 4      | 8      | 12    | 15    | 19     | 20     | 22.    | 24     | 27     | 30     |
| Individual    | mL     | mL     | mL    | mL    | mL     | mL     | mL     | mL     | mL     | mL.    |
| Cell 1        | 2329   | 3245   | 3588  | 3045  | -3181  | 1027   | 3322   | 3047   | 2851   | 2962   |
| Cell 2        | 2256   | 3145   | 3186  | 2904  | 3105   | 1001   | 3209   | 2950   | 2764   | 2861   |
| Cell 3        | 2293   | 3260   | 3315  | 3085  | 3198   | 1041   | 3280   | 3030   | 2836   | 2952   |
| Cell 4        | 2232   | 3076   | 3131  | 2937  | 3055   | 1003   | 3019   | 2902   | 2731   | 2840   |
| Accumulative  |        |        |       |       |        |        |        |        |        |        |
| Cell 1        | 2329   | 5575   | 9163  | 12208 | 15389  | 16416  | 19738  | 22785  | 25636  | 28593  |
| Cell 2        | 2256   | 5401   | 8587  | 11491 | 14596  | 15596  | 18806  | 21755  | 24519  | 27380  |
| Cell 3        | 2293   | 5552   | 8867  | 11952 | 15150  | 16191  | 19471  | 22501  | 25336  | 28283  |
| Cell 4        | 2232   | 5308   | 8440  | 11376 | 14432  | 15434  | 18453  | 21355  | 24086  | 26926  |

Table 19
Leachate Volumes Collected During the Soil Flushing Study for Phase II

| Sample Period | 1a     | 2     | 3     | 4     | 5      | 6      | 7      | 8      | 9      | 10     |
|---------------|--------|-------|-------|-------|--------|--------|--------|--------|--------|--------|
| Date          | 29-Apr | 2-May | 5-May | 8-May | 11-May | 14-May | 18-May | 21-May | 25-May | 28-May |
| Days          |        | 6     | 9 _   | 12    | 15     | 19     | 22     | 26     | 29     |        |
| Individual    | mL     | mL    | mL    | mL.   | mL     | mL     | mL     | mL     | mL     | mL     |
| Cell 1        | 3358   | 3090  | 1677  | 2952  | 3002   | 3016   | 3368   | 3023   | 3247   | 3042   |
| Cell 2        | 3283   | 2943  | 2830  | 2779  | 2826   | 2838   | 3270   | 2883   | 3185   | 2996   |
| Cell 3        | 3100   | 2999  | 2878  | 2837  | 2877   | 2892   | 3325   | 2892   | 3229   | 3037   |
| Cell 4        | 3208   | 2926  | 2814  | 2778  | 2803   | 2821   | 3197   | 2786   | 3159   | 2947   |
| Accumulative  |        |       |       |       |        |        | ·      |        |        |        |
| Days          |        | 3     | 6     | 9     | 12     | 15     | 19     | 22     | 26     | 29     |
| Cell 1        |        | 3090  | 4767  | 7719  | 10721  | 13737  | 17105  | 20128  | 23374  | 26416  |
| Cell 2        |        | 2943  | 5773  | 8552  | 11378  | 14216  | 17486  | 20369  | 23554  | 26550  |
| Cell 3        |        | 2999  | 5877  | 8713  | 11591  | 14483  | 17807  | 20699  | 23928  | 26965  |
| Cell 4        |        | 2926  | 5740  | 8518  | 11321  | 14142  | 17339  | 20125  | 23283  | 26230  |

a - denotes that leachate volume was not calculated into accumulative values

Table 20 Concentration of VOCs and SVOCs in Cell 1 Leachates

|                                |          | Phase I  |            | Phase II |          |              |  |
|--------------------------------|----------|----------|------------|----------|----------|--------------|--|
|                                | Period 1 | Period 2 | Composite  | Period 3 | Period 4 | Composite    |  |
| Chemical Analyte               | ug/L     | ug/L     | ug/L       | ug/L     | ug/L     | ug/L         |  |
| Volatile Organic Compounds     |          | ,-       |            |          |          |              |  |
| Benzene                        | 110      | 32       | 3          | -        | 16       | -            |  |
| Ethyl benzene                  | 1700     | -        | 4          | 1800     | 260      | •            |  |
| Toluene                        | 320      | 40       | . 5        | -        | 26       | <del>-</del> |  |
| Total xylenes                  | 7500     |          | -          | -        | •        | •            |  |
| Semivolatile Organic Compounds |          |          |            | •        |          |              |  |
| Butylbenzylphthalate           | -        | -        | -          | 540      | -        | •            |  |
| Bis(2-ethylhexyl)phthalate     | -        | •        |            | 61000    | 3200     | 6500         |  |
| Di-n-butylphthalate            | 20       | -        | -          | -        | -        | -            |  |
| Diethyl phthalate              | 86       | -        | -          | -        | -        | -            |  |
| Dimethyl phthalate             | 26       | •        | <b>-</b> ; | -        | -        | <b>-</b> .   |  |
| Di-n-octyl phthalate           | -        | -        | -          | 190      | •        | 24           |  |
| Total phthalates (additional)  |          | •        | •          | -        |          | · -          |  |
| LEACHATE VOLUME (L)            | 2.329    | 2.962    | 28.598     | 3.09     | 3.042    | 26.416       |  |

Table 21
Concentration of VOCs and SVOCs in Cell 2 Leachates

|                                |          | Phase I    |            | Phase II |          |           |  |  |
|--------------------------------|----------|------------|------------|----------|----------|-----------|--|--|
|                                | Period 1 | Period 2   | Composite  | Period 3 | Period 4 | Composite |  |  |
| Chemical Analyte               | ug/L     | ug/L       | ug/L       | ug/L     | ug/L     | ug/L      |  |  |
| Volatile Organic Compounds     |          |            |            |          |          |           |  |  |
| Benzene                        | 120      | 4.         | <b>3</b>   | 7        | 6        | 6         |  |  |
| Ethyl benzene                  | 71       | -          | 4          | -        | -        | 3         |  |  |
| Toluene                        | _ 310    | 5:_        | <b>5</b>   | 18       | : 1:6°   | .14       |  |  |
| Total xylenes                  | 840      | •          |            |          | -        | <b>~</b>  |  |  |
| Semivolatile Organic Compounds |          |            |            |          |          |           |  |  |
| Butylbenzylphthalate           | -        | -          | -          | <u>.</u> | -        | -         |  |  |
| Bis(2-ethylhexyl)phthalate     | -        | •          |            | :=:      | -        | -         |  |  |
| Di-n-butylphthalate            | -        | <b>-</b> . | -          | ■,       | -        | -         |  |  |
| Diethyl phthalate              | 11       | -          | -          | -        | 2        | •         |  |  |
| Dimethyl phthalate             | -        | •          | -          | -        | -        | -         |  |  |
| Di-n-octyl phthalate           | -        | -          |            | -        | -        | -         |  |  |
| Total phthalates (additional)  |          | -          | <u>.</u> . | -        | •        | -         |  |  |
| LEACHATE VOLUME (L)            | 2.256    | 2.861      | 27.38      | 3.283    | 2.996    | 26.55     |  |  |

Table 22
Concentration of VOCs and SVOCs in Cell 3 Leachates

|                                |          | Phase I      |            | Phase II   |          |           |  |
|--------------------------------|----------|--------------|------------|------------|----------|-----------|--|
|                                | Period 1 | Period 2     | Composite  | Period 3   | Period 4 | Composite |  |
| Chemical Analyte               | ug/L     | ug/L         | ug/L       | ug/L       | ug/L     | ug/L      |  |
| Volatile Organic Compounds     |          | ·            |            |            |          |           |  |
| Benzene                        | 48       | 3            | <b>-</b> 3 | -          | 2        | -         |  |
| Ethyl benzene                  | 4        | -            | 2          |            | 3        | -         |  |
| Toluene                        | 100      | <b>3</b> . ; | 5          | .= .       | 3        | -         |  |
| Total xylenes                  | 410      | -            | •          | <b>-</b> . | ,=       | -         |  |
| Semivolatile Organic Compounds |          |              |            |            |          |           |  |
| Butylbenzylphthalate           | -        | -            | 2          | •          | -        | 6         |  |
| Bis(2-ethylhexyl)phthalate     | -        | -            | -          | -          | -        | -         |  |
| Di-n-butylphthalate            | -        | -            | -          | -          | -        | -         |  |
| Diethyl phthalate              | 12       | -            | .=         | -          | -        | -         |  |
| Dimethyl phthalate             | -        |              | -          | -          |          | -         |  |
| Di-n-octyl phthalate           | -        | -            | -          | -          | -        | -         |  |
| Total phthalates (additional)  | -        |              | -          | •          | -        | -         |  |
| LEACHATE VOLUME (L)            | 2.293    | 2.952        | 28.288     | 3.1        | 3.037    | 26.965    |  |

Table 23
Concentration of VOCs and SVOCs in Cell 4 Leachates

|                                |          | Phase I  |           | Phase II |              |            |  |
|--------------------------------|----------|----------|-----------|----------|--------------|------------|--|
|                                | Period 1 | Period 2 | Composite | Period 3 | Period 4     | Composite  |  |
| Chemical Analyter              | ug/L     | ug/L     | ug/L      | ug/L     | ug/L         | ug/L       |  |
| Volatile Organic Compounds     |          |          |           |          |              |            |  |
| Benzene                        | 10       | . •      | -         | •        | 15           | 13         |  |
| Ethyl benzene                  | · - '    | -        | 1         | . •      | . =          |            |  |
| Toluene                        | 2        | •        | 2         | -        | 21           | 20         |  |
| Total xylenes                  | 64       | •        | •         | -        | -            | . •        |  |
| Semivolatile Organic Compounds |          |          |           |          |              | -          |  |
| Butylbenzylphthalate           | -        | -        | -         | 190      | -            | -          |  |
| Bis(2-ethylhexyl)phthalate     | -        | •        | -         | 490      | 360          | 38         |  |
| Di-n-butylphthalate            | -        | •        | -         | .=.      | -            |            |  |
| Diethyl phthalate              | 2        | -        | -         | •        | : <b>.</b> : | -          |  |
| Dimethyl phthalate             | -        | -        | •         |          | -            | -          |  |
| Di-n-octyl phthalate           | -        | -        | -         | -        | -            | -          |  |
| Total phthalates (additional)  | •        | <u>-</u> | -         | -        | •            | <u>-</u> . |  |
| LEACHATE VOLUME (L)            | 2.232    | 2.84     | 26,926    | 3.208    | 2.947        | 26.23      |  |

Table 24
Concentration of Analytes in Final Soil Samples
Used in the Soil Flushing Study

| Cell Number                              | Cell 1         | Cell 2       | Cell 3   | Cell 4   |
|--|----------------|--------------|----------|----------|
| Sample Identification                    | ST-2A          | ST-3A        | ST-3A    | ST-4D    |
| Sample Depth (inches)                    | 72 to 96       | 24 to 48     | 24 to 48 | 12 to 36 |
| Chemical Analyte                         | ug/kg          | ug/kg        | ug/kg    | ug/kg    |
| Volatile Organic Compounds               |                |              |          |          |
| Ethyl benzene                            | 29000          | -            | -        | 9        |
| 4-methyloctane                           | ` <del></del>  | -            | -        | -        |
| Total xylenes                            | -              | -            | -        | -        |
| 2,5-dimethylnonane                       | <del>.</del>   | . •          | -        | -        |
| Propylcyclohexane                        | : <del>_</del> | -            | -        | -        |
| Ethylmethylbenzene                       | -              | •            | -        | -        |
| 2-methylnonane                           | -              | -            | -        | -        |
| Total cyclohexanes                       | e              | •            | • .      | •        |
| Semivolatile Organic Compounds           |                |              |          |          |
| Butylbenzylphthalate                     | 1800           | •            | •        | 600      |
| Bis(2-ethylhexyl)phthalate               | 110000         | 20000        | 11000    | 310000   |
| Di-n-butylphthalate                      | 600            | •            | -        | -        |
| Diethyl phthalate                        | -              | -            | -        | •.       |
| Dimethyl phthalate                       |                | -            | -        | •        |
| Di-n-octyl phthalate                     | 2500           | 310          | 440      | 1100     |
| Total phthalates (additional)            |                | •            | -        | •        |
| Anthracene                               | <del>.</del>   | -            | -        | •        |
| Benzo(a)anthracene                       | -              | -            | -        | •        |
| Benzo(b)fluoranthene                     | •              | <del>.</del> | -        | -        |
| Benzo(k)fluoranthene                     | -              | -            | -        | •        |
| Benzo(a)pyrene                           | •              | •            | 120      | ÷        |
| Chrysene                                 | •              | -            | -        | •        |
| Fluoranthene                             | •              | •            | •        | -        |
| Indeno(1,2,3-cd)pyrene                   | -              | -            | -        | •        |
| Naphthalene                              |                | -            | -        | -        |
| Phenanthrene                             | •              | -            | -        | -        |
| Pyrene                                   | •              | -            | •        | •        |
| Dibenz(a,h)anthracene                    | •              | -            | •        | .=       |
| 4-Hydroxy-4-methyl-2-pentanone           | •              | •            | -        | •        |
| Hexadecanoic acid                        | ; <del>*</del> | -            | -        | -        |
| 2-Ethylhexyl diphenyl ester phosphoric a | •              | -            | -        | -        |
| 2-Methylpropyl ester octadecanoic acid   | <b>1</b>       | <b>=</b>     | -        | •        |
| Total alkanes                            |                | -            | -        | -        |

Table 25
Total Amount of VOCs and SVOCs in Cell 1 Leachates

|                                |              | Phase I      |           |            | Phase II |           | Phase I | Phase II     |
|--------------------------------|--------------|--------------|-----------|------------|----------|-----------|---------|--------------|
|                                | Period 1     | Period 2     | Composite | Period 3   | Period 4 | Composite | SUM     | SUM          |
| Chemical Analyte               | ug           | ug           | ug        | ug         | ug       | ug        | ug      | ug           |
| Volatile Organic Compounds     |              |              |           |            |          |           |         |              |
| Benzene                        | -256         | 95           | - 86      |            | 49       | •         | 437     | 49           |
| Ethyl benzene                  | 3959         | -            | ·114      | 5562       | 791      | •         | 4074    | 6353         |
| Toluene                        | 745          | 118          | 143       |            | 79       | -         | 1007    | <b>79</b>    |
| Total xylenes                  | 17468        | <del>.</del> |           | . <b>-</b> | . •      |           | 17468   | . 0          |
| Semivolatile Organic Compounds |              |              |           | ·          | •        |           |         |              |
| Butylbenzylphthalate           | -            | -            | -         | 1669       | •        | -         | 0       | 1669         |
| Bis(2-ethylhexyl)phthalate     | -            | -            | -         | 188490     | 9734     | 171704    | 0       | 369928       |
| Di-n-butylphthalate            | 47           | -            | -         | -          | -        | -         | 47      | 0            |
| Diethyl phthalate              | 200          | -            | -         | -          | -        | -         | 200     | 0            |
| Dimethyl phthalate             | 61           | •            | -         | -          | -        | -         | 61      | 0            |
| Di-n-octyl phthalate           | -            | -            | -         | 587        | •        | 634       | 0       | <b>12</b> 21 |
| Total phthalates (additional)  | , <b>-</b> - | -            | -         | -          | -        | -         | 0       | 0            |

Table 26
Total Amount of VOCs and SVOCs in Cell 2 Leachates

|                               |          | Phase I  |           | Phase II  |          |           | Phase I | Phase II |
|-------------------------------|----------|----------|-----------|-----------|----------|-----------|---------|----------|
|                               | Period 1 | Period 2 | Composite | Period 3  | Period 4 | Composite | SUM     | SUM      |
| CHEMICAL ANALYTE              | ug       | ug       | ug        | ug        | ug       | ug        | ug      | ug       |
| Volatile Organic Compounds    |          |          |           |           |          |           |         |          |
| Benzene                       | 271      | 11       | 82        | <b>23</b> | 18       | 159       | 364     | 200      |
| Ethyl benzene                 | 160      |          | 110       | •         | •        | 80        | 270     | 80       |
| Foluene Foluene               | 699      | 14       | 137       | 59        | 48       | 372       | 850     | 479      |
| Total xylenes                 | 1895     | ·<br>•   | -         | -         |          | •         | 1895    | 0        |
| Semivolatile Organic Compou   | ınds     |          |           |           | •        |           |         |          |
| Butylbenzylphthalate          | •        | -        | -         | -         | -        | •.        | 0       | 0        |
| Bis(2-ethylhexyl)phthalate    | -        | -        | -         | •.        | -        | •         | 0       | 0        |
| Di-n-butylphthalate           | -        | -        | -         | -         | -        | -         | 0       | 0        |
| Diethyl phthalate             | 25       | •        | -         | -         | 6        | •         | 25      | 6        |
| Dimethyl phthalate            | •        |          |           | -         | -        | -         | 0       | 0        |
| Di-n-octyl phthalate          | -        | -        | -         | -         | •        | -         | 0       | 0        |
| Fotal phthalates (additional) | •        | -        | -         | -         | -        | -         | 0       | 0        |

Table 27
Total Amount of VOCs and SVOCs in Cell 3 Leachates

|                                | Phase I  |          |                                       | Phase II |                      |             | Phase I | Phase II |
|--------------------------------|----------|----------|---------------------------------------|----------|----------------------|-------------|---------|----------|
|                                | Period 1 | Period 2 | Composite                             | Period 3 | Period 4             | Composite   | SUM     | SUM      |
| CHEMICAL ANALYTE               | ug       | ug       | ug                                    | ug       | ug                   | ug          | ug      | ug       |
| Volatile Organic Compounds     |          |          |                                       |          |                      | <u> </u>    |         |          |
| Benzene                        | 110      | 9        | 85                                    | -        | 6                    | -           | 204     | 6        |
| Ethyl benzene                  | 9        | -        | 57                                    | -        | 9                    | -           | 66      | 9        |
| Toluene                        | 229      | 9        | 141                                   | -        | 9                    | •           | .380    | 9        |
| Total xylenes                  | 940      |          | -                                     |          | -                    | -           | 940     | . 0      |
| 14.14.1                        |          | *=       | · · · · · · · · · · · · · · · · · · · |          |                      |             |         |          |
| Semivolatile Organic Compounds |          |          |                                       |          | •                    |             |         |          |
| Butylbenzylphthalate           | •        | -        | 57                                    | - · .    | . <b>=</b> , . , , , | <b> 162</b> | 57      | 162      |
| Bis(2-ethylhexyl)phthalate     | -        | •        | -                                     | .=       | -                    | -           | 0       | 0        |
| Di-n-butylphthalate            | -        | •        | -                                     | -        | -                    | •           | 0       | 0        |
| Diethyl phthalate              | 28       | -        | -                                     |          | :=                   | -           | 28      | 0        |
| Dimethyl phthalate             | -        | -        | -                                     | -        | -                    | -           | 0       | 0        |
| Di-n-octyl phthalate           | -        | -        | -                                     |          | -                    | -           | 0       | 0.       |
| Total phthalates (additional)  | -        | •        | -                                     | -        | -                    | -           | 0       | 0        |

Table 28

Total Amount of VOCs and SVOCs in Cell 4 Leachates

|                                | Phase I    |          |           | Phase II |          |            | Phase I                                 | Phase II      |
|--------------------------------|------------|----------|-----------|----------|----------|------------|---|---------------|
|                                | Period 1   | Period 2 | Composite | Period 3 | Period 4 | Composite  | SUM                                     | SUM           |
| Chemical Analyte               | ug         | ug       | ug        | ug       | ug       | ug         | ug                                      | ug            |
| Volatile Organic Compounds     |            |          |           |          |          |            | , , , , , , , , , , , , , , , , , , , _ |               |
| Benzene                        | 22         | -        | -         | -        | 44       | 341        | 22                                      | 385           |
| Ethyl benzene                  | -          | -        | 27        | -        | _        | -          | 27                                      | 0             |
| Toluene                        | 4          | -        | 54        | -        | 62       | 525        | 58                                      | 586           |
| Total xylenes                  | 143        | -        | -         | -        | -        | -          | 143                                     | 0             |
| Semivolatile Organic Compounds |            |          |           |          |          |            |   | 4             |
| Butylbenzylphthalate           |            | -        | - i       | 610      | 5.552    | • • •      | 0                                       | - 610         |
| Bis(2-ethylhexyl)phthalate     | <b>1-</b>  | -        | -         | 1572     | 1061     | 997        | 0                                       | 3630          |
| Di-n-butylphthalate            | · · · · == | -        | •         | -        | •        |            | 0                                       | - · · · · · O |
| Diethyl phthalate              | 4          | -        | -         | -        | -        | -          | 4                                       | 0             |
| Dimethyl phthalate             | -          | -        | -         | -        |          | •          | 0                                       | . 0           |
| Di-n-octyl phthalate           | -          | -        | -         | -        | -        | • .        | 0                                       | 0             |
| Total phthalates (additional)  | -          | -        | -         | -        | -        | . <b>-</b> | 0                                       | 0             |

<sup>\*</sup> denotes analyte was found in flushing solution

Table 29

Total Percent and Amount of Contaminants Removed From Soils During Flushing

|                               | Cell 1    |           |          |       |                                       | Cell 2   |              |      |  |
|-------------------------------|-----------|-----------|----------|-------|---------------------------------------|----------|--------------|------|--|
|                               | Phase I   | Phase II  | Tenax    | Soil  | Phase I                               | Phase II | Tenax        | Soil |  |
| <b>Chemical Analyte</b>       | %\ug      | %\ug      | %\ug     | ug    | %\ug                                  | %\ug     | %\ug         | ug   |  |
| VOCs                          |           |           |          |       |                                       |          | <u> </u>     |      |  |
| Benzene                       | 81\437    | 9\49      | 10\54    | 0     | 52\364                                | 29\200   | 19\130       | 0    |  |
| Ethyl benzene                 | 16\4074   | 24\6353   | 2\430    | 15083 | 66\270                                | 19\80    | 15\62        | 0    |  |
| Toluene                       | 82\1007   | 6\79      | 11\140   | 0     | 50\850                                | 28\479   | 21\360       | 0    |  |
| Total xylenes                 | 100\17468 | 0/0       | 0\0      | 0     | 100\1895                              | 0/0      | 0/0          | 0    |  |
| SVOCs                         |           |           |          |       | · · · · · · · · · · · · · · · · · · · | - ,      |              | *    |  |
| Butylbenzylphthalate          | 0/0       | 64\1669   | -        | 936   | 0/0                                   | 0/0      | . <b>-</b> _ | 0    |  |
| Bis(2-ethylhexyl)phthalate    | 0/0       | 87\369928 | <u>-</u> | 57211 | 0\0                                   | 0\0      | -            | 7176 |  |
| Di-n-butylphthalate           | 13\47     | 0/0       | -        | 312   | 0/0                                   | 0/0      | -            | 0    |  |
| Diethyl phthalate             | 100\200   | 0/0       | -        | 0     | 81\25                                 | 19\6     | -            | 0    |  |
| Dimethyl phthalate            | 100\61    | 0\0       | •        | 0     | 0/0                                   | 0/0      | -            | 0    |  |
| Di-n-octyl phthalate          | 0\0       | 48\1221   | -        | 1300  | 0/0                                   | 0/0      | -            | 111  |  |
| Total phthalates (additional) | 0/0       | 0/0       | -        | 0     | 0/0                                   | 0/0      | -            | 0    |  |

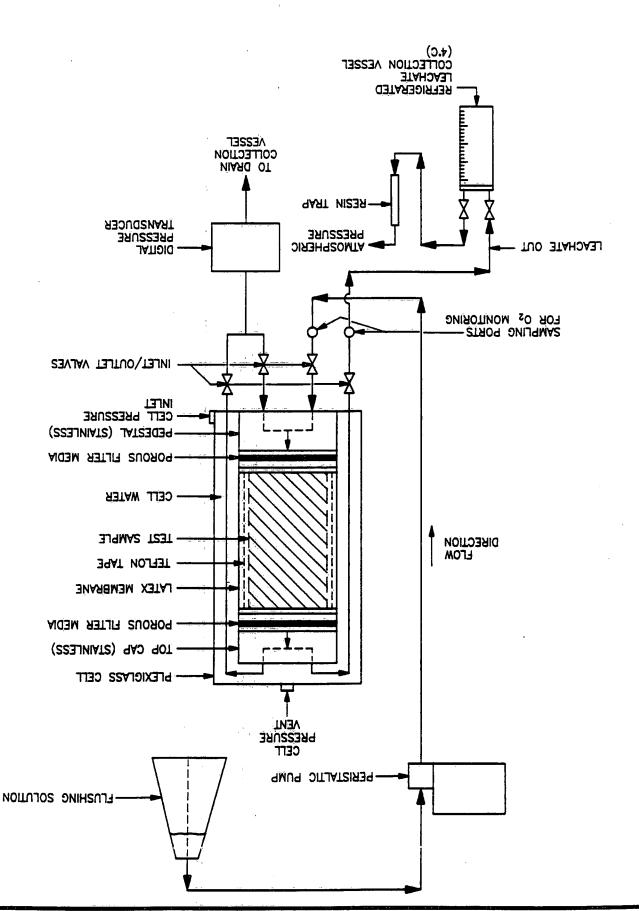
Table 29 (cont.)

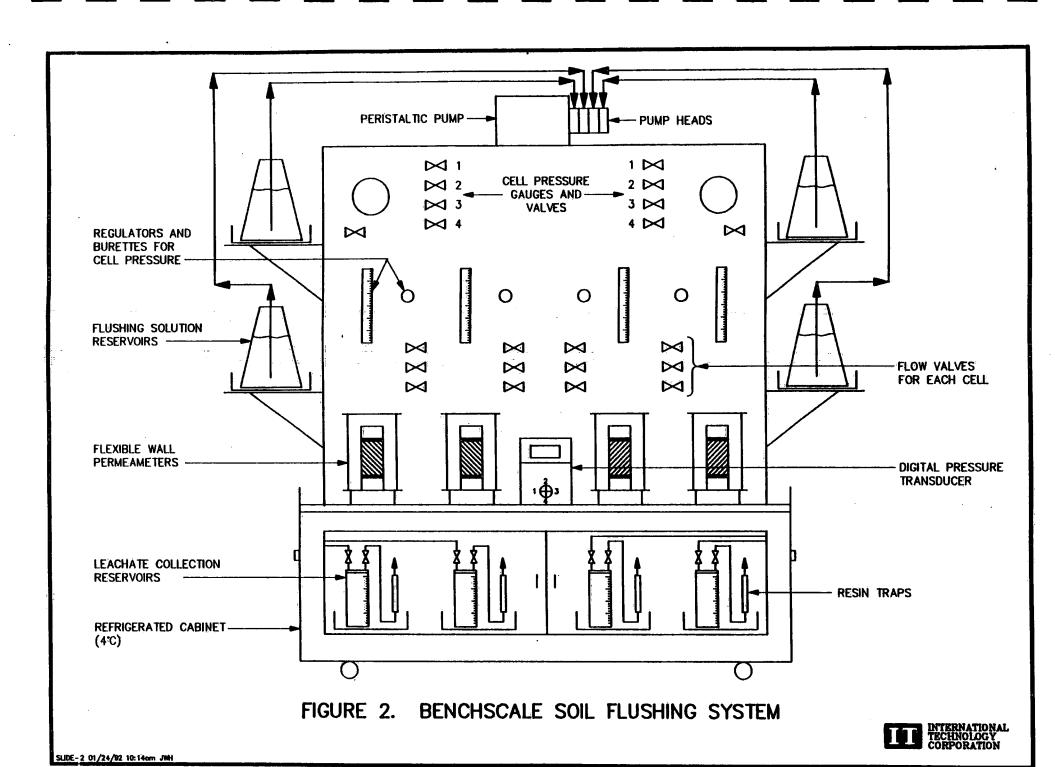
Total Percent and Amount of Contaminants Removed From Soils During Flushing

|                               | Cell 3  |          |        |      |         | Cell 4   |          |        |  |
|-------------------------------|---------|----------|--------|------|---------|----------|----------|--------|--|
|                               | Phase I | Phase II | Tenax  | Soil | Phase I | Phase II | Tenax    | Soil   |  |
| Chemical Analyte              | %\ug    | %\ug     | %\ug   | ug   | %\ug    | %\ug     | %\ug     | ug     |  |
| VOCs                          |         |          |        |      |         |          |          |        |  |
| Benzene                       | 71\204  | 2\6      | 27\79  | 0    | 5\22    | 93\385   | 1\6      | 1      |  |
| Ethyl benzene                 | 63\66   | 9\9      | 29\30  | 0    | 73\27   | :0/0     | 27\10    | 0      |  |
| Toluene                       | 60\380  | 1\9      | 38\240 | - 0  | 9\58    | 89\586   | 2\11     | 1      |  |
| Total xylenes                 | 100\940 | 0/0      | 0/0    | 0    | 100\143 | 0/0      | 0//0     | 0      |  |
| SVOCs                         |         |          |        |      | •       |          | •        |        |  |
| Butylbenzylphthalate          | 26\57   | 74\162   | -      | Ó    | 0\0     | 66\610   | <u>-</u> | 321    |  |
| Bis(2-ethylhexyl)phthalate    | 0/0     | 0/0      | -      | 5407 | 0/0     | 2\3630   |          | 165990 |  |
| Di-n-butylphthalate           | 0/0     | 0/0      | -      | 0    | 0/0     | 0\0      | -        | 0      |  |
| Diethyl phthalate             | 100\28  | 0/0      | -      | 0    | 100\4   | 0/0      | -        | 0      |  |
| Dimethyl phthalate            | 0/0     | 0/0      | -      | Ō    | 0\0     | 0\0      | -        | 0      |  |
| Di-n-octyl phthalate          | 0/0     | 0/0      | _      | 216  | 0/0     | 0/0      | -        | 589    |  |
| Total phthalates (additional) | 0/0     | 0/0      | •      | 0    | 0/0     | 0\0      | -        | 0      |  |



## FIGURE 1. FLEXIBLE WALL PERMEAMETER





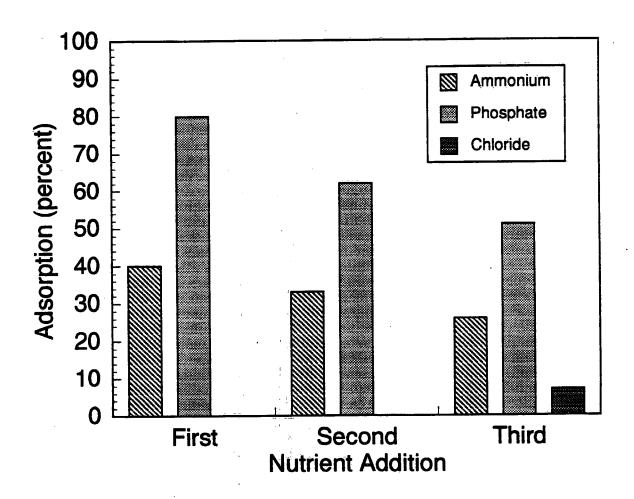


Figure 3

Nutrient adsorption. Nutrients were added to a 9 parts water and 1 part soil slurry as three spikes with measurement of adsorption performed after 24 hours of incubation following each spike.

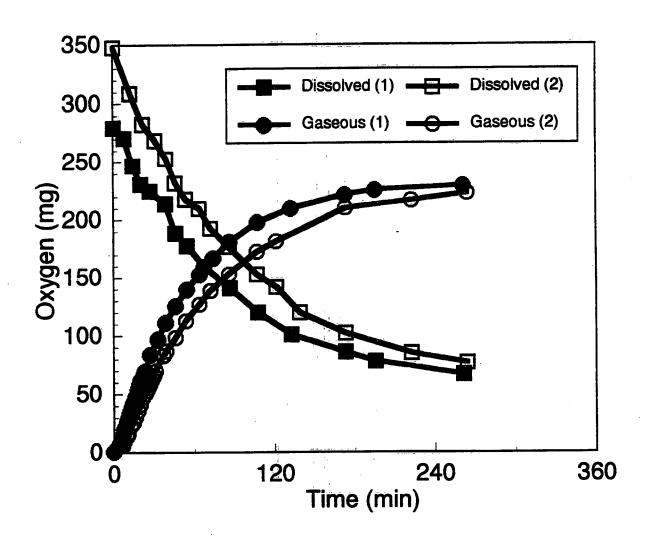


Figure 4

Oxygen delivery potential. The decomposition of hydrogen peroxide into gaseous oxygen and the corresponding residual total soluble oxygen including dissolved oxygen and hydrogen peroxide was measured.

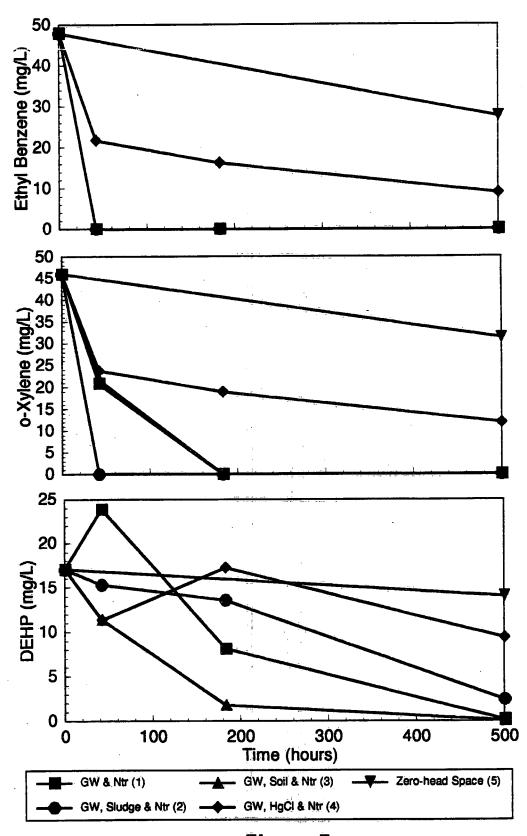
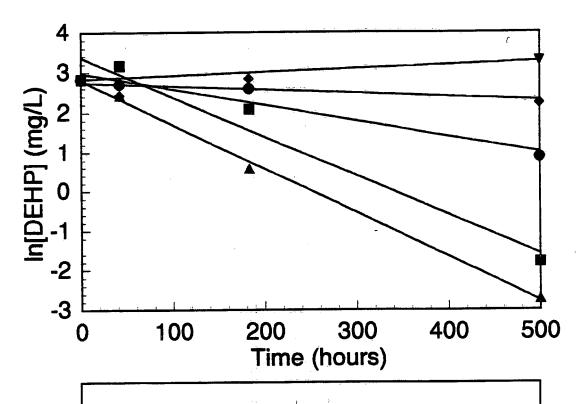


Figure 5

Ethyl Benzene, xylene, and DEHP removal from groundwater batch treatability tests.
L. E. Carpenter, IT Project No. 408474



- **GW & Ntr:**  $k = -9.898E-3 \text{ hrs}^{-1}$ ,  $t_{1/2} = 70 \text{ hours}$
- GW, Sludge & Ntr:  $k = -3.933E-3 \text{ hrs}^{-1}$ ,  $t_{1/2} = 176 \text{ hours}$
- **△** GW, Soil & Ntr: k = -1.115E-2 hrs<sup>-1</sup>,  $t_{1/2} = 62$  hours
- GW, HgCl & Ntr:  $k = -8.668E-4 \text{ hrs}^{-1}$ ,  $t_{1/2} = 800 \text{ hours}$
- ▼ Zero Head Space: infinite

Figure 6
Biodegradation kinetic constants and half lives for DEHP in each batch treatment.

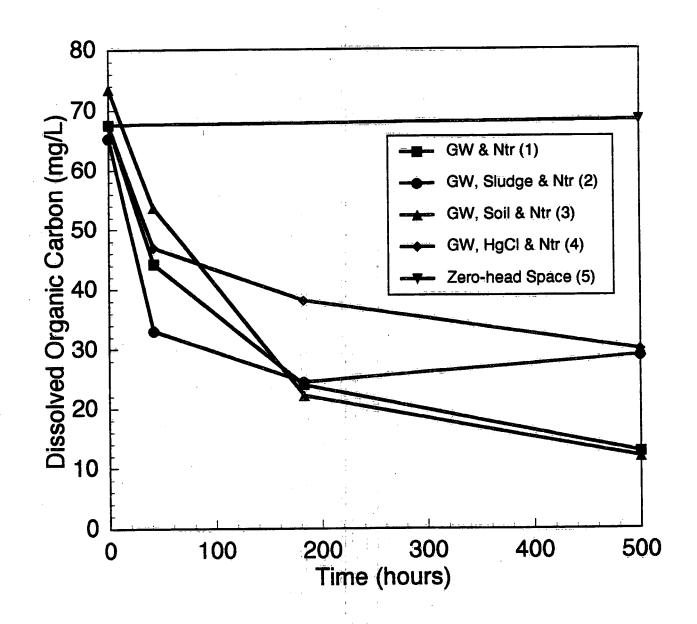


Figure 7

Mean dissolved organic carbon content of groundwater during batch treatment.

L. E. Carpenter and Co. IT Project No. 408474.

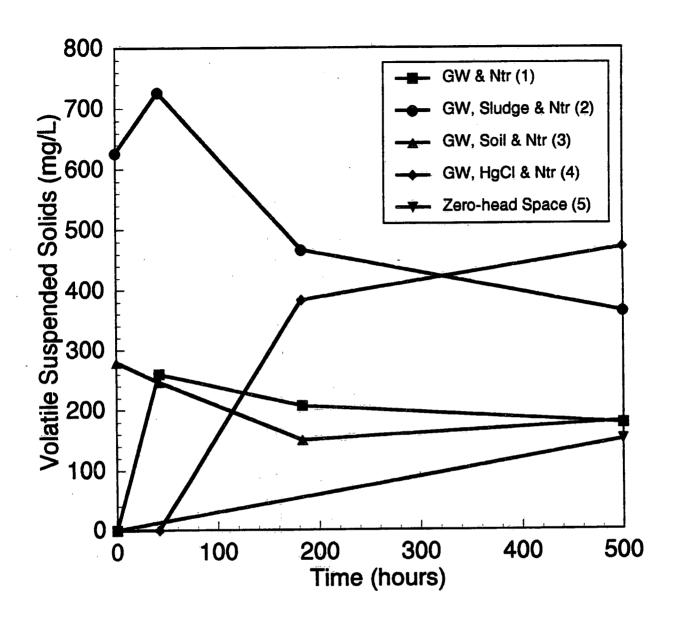


Figure 8

Mean volatile suspended solids analysis of groundwater treatments.

L. E. Carpenter and Co. IT Project No. 408474.

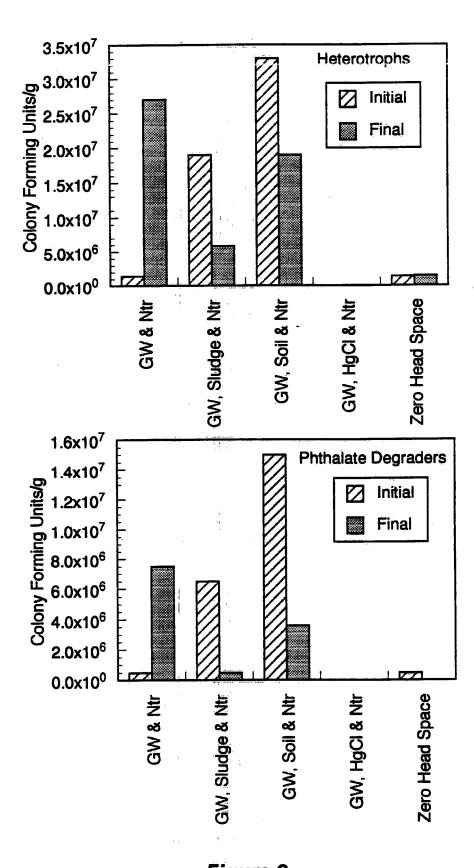


Figure 9
Initial and final microbial population density in groundwater batch treatability tests

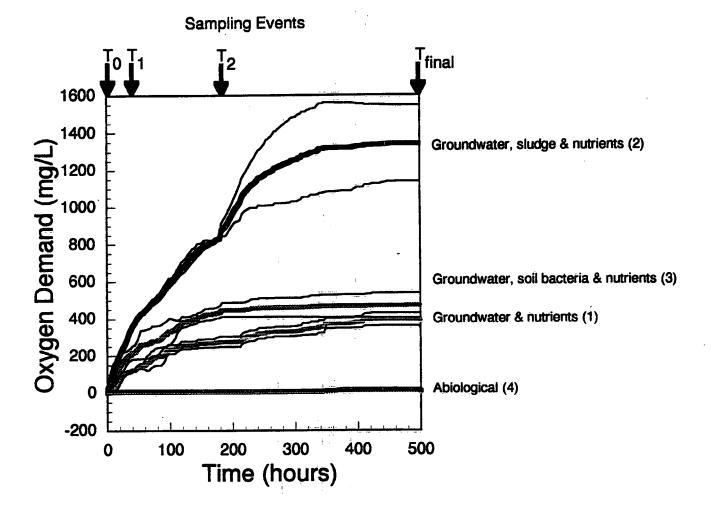
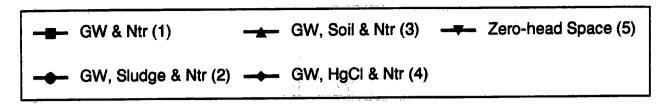


Figure 10

Microbial respiration in groundwater batch treatability studies. Thick lines present the mean of three values for each treatment. Thin lines indicate  $\pm$  the standard deviation for each data point. Data collection frequency was 2 hours.



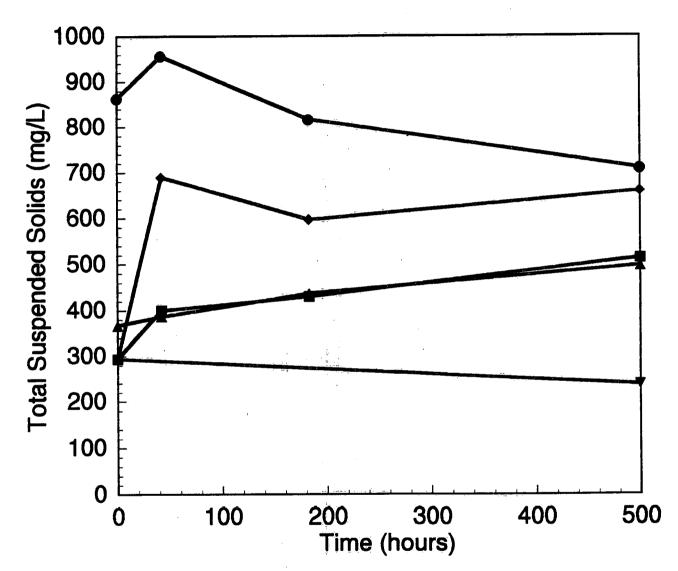


Figure 11

Mean total suspended solids analysis of groundwater treatments.

L. E. Carpenter and Co. IT Project No. 408474.

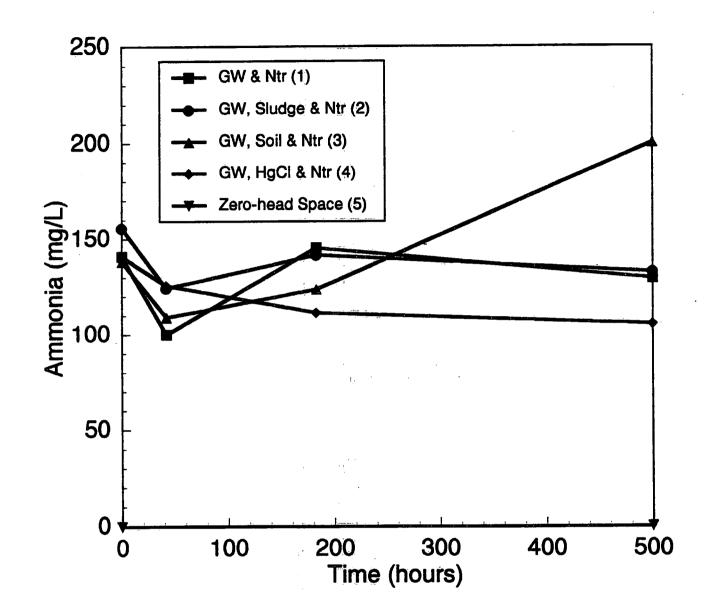


Figure 12

Ammonia utilization during groundwater respiration evaluation.

L. E. Carpenter and Co. IT Project No. 408474.

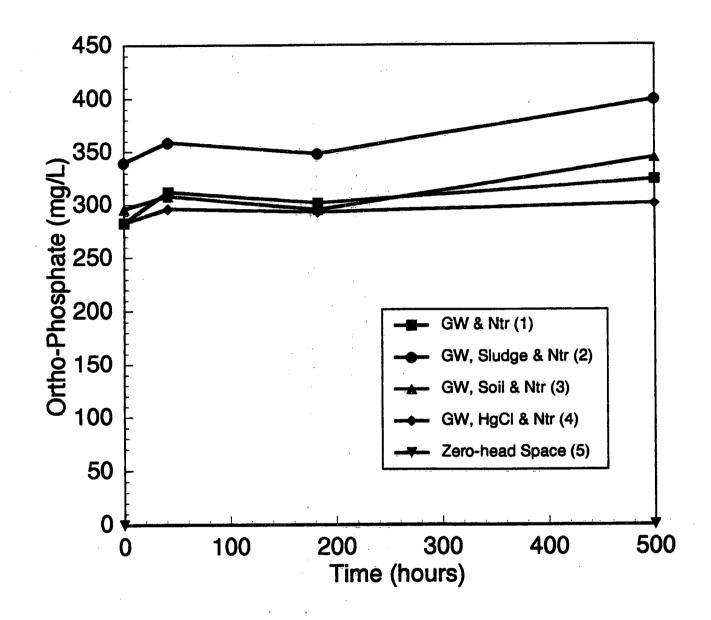


Figure 13

Phosphate utilization during groundwater batch study.

L. E. Carpenter and Co. IT Project No. 408474.

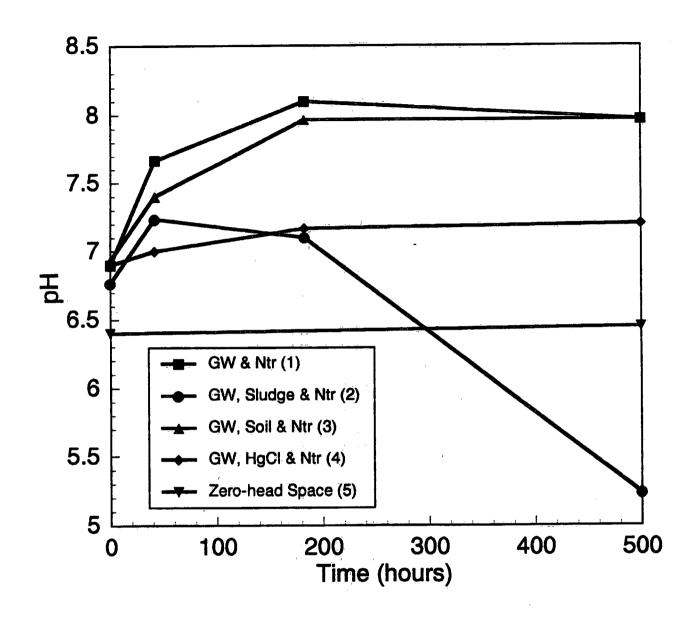


Figure 14

Change in pH during groundwater batch treatment.
L. E. Carpenter and Co. IT Project No. 408474.

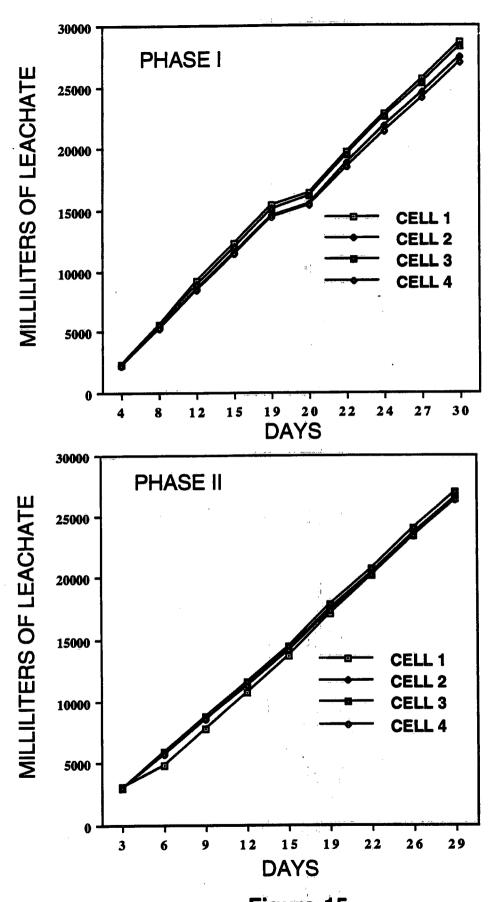


Figure 15
Milliliters of leachate collected for four samples during the two phases of the soil flushing study



# BIOREMEDIATION AND SOIL FLUSHING TREATABILITY STUDY REPORT L. E. CARPENTER AND CO. PART 2 OF 2

APPENDICES

PREPARED FOR L. E. CARPENTER AND CO.

PREPARED BY

JUNE 1992 PROJECT NO. 408474

Regional Office 312 Directors Drive • Knoxville, Tennessee 37923 • 615-690-3211

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# Appendix A

Sampling Plan

### 1.0 <u>INTRODUCTION</u>

Roy F. Weston, Inc. (WESTON) has prepared this document to provide documentation on the locations from which the L.E Carpenter treatability study samples will be collected. Furthermore, this document will discuss the rationale used to determine the selection of these locations.

### 2.0 SAMPLE LOCATION SELECTION CRITERIA

Soil sampling point locations were selected using two primary criteria: location with respect to known sampling points (i.e., test pit samples, hand auger samples, and monitoring well locations), and concentration of contaminants of concern detected in those previous samples. For the former criteria, locations were first selected based upon their proximity to previous sampling locations with documented contaminant concentration. Specifically, soil sampling locations were chosen which were indicative of low, moderate, and high ranges of contaminants of concern (mainly bis(2-ethylhexyl)phthalate (DEHP)). Furthermore, soil sample locations were selected which should be indicative of general metals concentrations on-site, and not metals "hot spot" conditions.

Wells were also selected using two primary criteria; location of the well with respect to known organic compound sources and known concentrations of volatile organic compounds (VOCs) and semi-volatile organic compounds (SVOCs) in the wells. For the former criteria, wells were selected from upgradient, source area and downgradient locations. For the latter criteria, wells which showed low, intermediate and high VOC and SVOC concentrations were selected.

### 3.0 SELECTED LOCATIONS: SOIL

Figure 3-1 indicates the selected locations for collection of background, low, medium, and high concentration Shelby tube samples, in relation with salient site features, such as monitoring wells, property boundaries and the Rockaway River. This map has been compared with other site maps which depict the location of test pit samples which were collected as part of the Remedial Investigation.

The background sample (indicated on Figure 3-1 by S.B) will be collected from a location 100' south of monitoring well MW-1, approximately 60' from the former production well. This is the location from which test pit sample TP78 was collected. This sample indicated minimal concentrations of DEHP (260 ug/kg, also detected in the laboratory method blank), and no volatile organic contamination. This area should also be indicative of the fill material (mine tailings) historically placed at the site.

The low-concentration sample (indicated on Figure 3-1 by S,L) will be located approximately

125' due north of monitoring well cluster MW-11, approximately 60' from recovery well RW-3. This is in the area of test pit sample TP74 which indicated 2089 ug/kg of total base neutral and acid extractable compounds. This location is outside the calculated area of influence resulting from the floating product lens associated with monitoring well MW-11s.

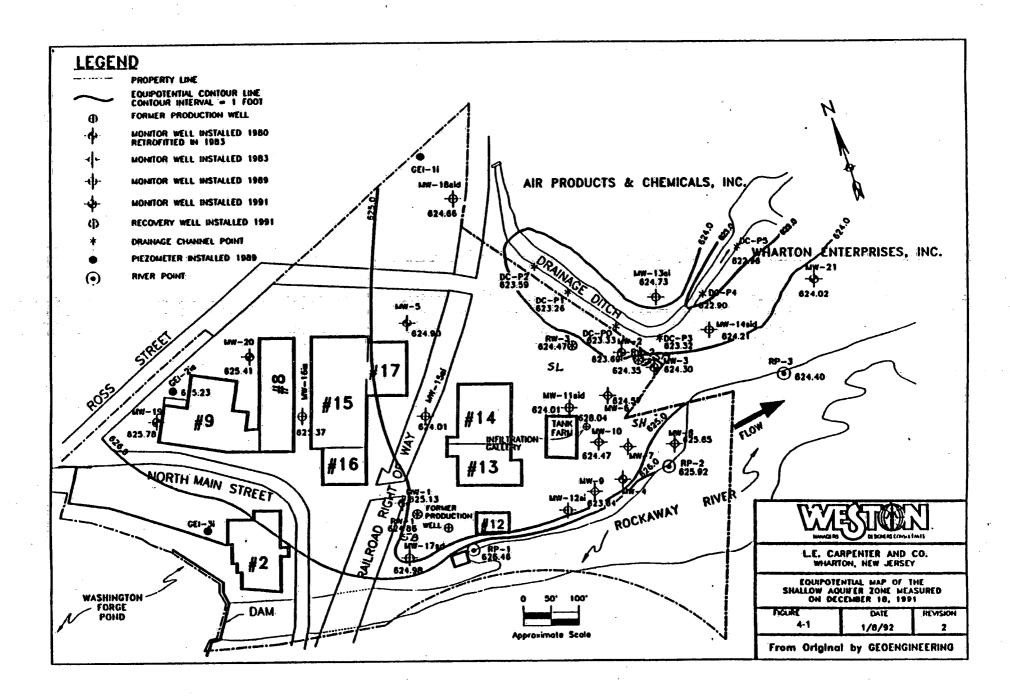
The medium concentration Shelby tube sample will be collected from the location on Wharton Enterprises property, approximately 90' from monitoring well cluster MW-14, adjacent to monitoring well MW3. This location is depicted on Figure 3-1 by S<sub>2</sub>M, and is in the vicinity of test pits TP3 and TP85, which indicated semivolatile organics in the range of 109,000 ug/kg to 510,000 ug/kg. Additionally, volatile organics (primarily xylenes) were detected at a total concentration of 45,800 ug/kg in TP3 at a depth of 4.5'-5', while PCBs were detected at 14 ppm in the surface soil (0'-.5') in TP3.

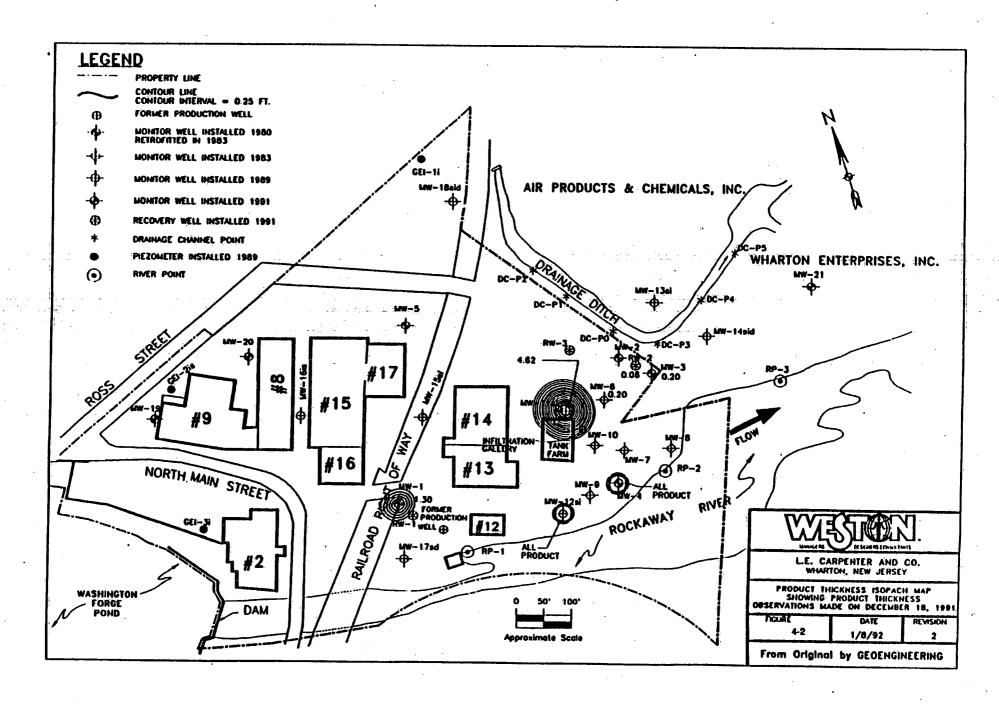
The high concentration Shelby tube samples will be collected from a location (as depicted on Figure 3-1 by S<sub>t</sub>H) equidistant from monitoring wells MW6 and MW8, adjacent to property line which L.E. Carpenter shares with Wharton Enterprises. This location was bounded by test pit samples TP5, TP6, TP7, and TP8, which ranged in concentration from 3.1 x 10<sup>6</sup> ug/kg to 10 x 10<sup>6</sup> ug/kg DEHP and 200 ug/kg to 532,000 ug/kg total targeted volatiles. If sample collection from this location should prove impracticable in the field, an alternate location (as depicted on Figure 3-1 by AS<sub>t</sub>H), approximately 25' west of the northwest corner of the tank farm containment pad, will be used. Test pit samples (TP42-TP48) collected in this area indicated DEHP concentrations ranging from 7.3 x 10<sup>6</sup> to 30 x 10<sup>6</sup> ug/kg.

In all cases, the Shelby tube samples will be collected in two adjacent 30" lengths (60" total length) immediately above the water table. Since the contaminants of concern have the capability of forming a lighter than water, non-aqueous phase layer (LNAPL), the highest concentrations in vadose zone soils would be expected between the water table and the historic high water table. These soils would be the primary focus of an in-situ treatment scheme. Therefore, soil samples collected from this depth would most closely delineate actual in-situ treatment parameters.

## 4.0 SELECTED LOCATIONS - GROUNDWATER

Figure 4-1 shows the equipotential surface of the shallow aquifer zone measured in November of 1991. This figure illustrates the main aspects of the shallow hydrogeology of this site. In the vicinity of the site, the Rockaway River is a recharge boundary. Shallow aquifer zone groundwater flow vectors along the river area oriented to the north. Toward the center of the site, these vectors swing back to the east and are oriented parallel to the Rockaway River. Figure 4-2 shows that there are two main source areas; a localized area around MW-1 and a larger area around the former impoundment area east of the tank farm. The western portion of the treatability study area is considered to be upgradient and the eastern portion is considered to be downgradient. Three general well locations; upgradient, source area and downgradient; can be selected using this information. With





respect to the former impoundment area, MW-15S can be considered an upgradient well. Source area wells include MW-3, MW-6, and MW-11S. MW-14S is the downgradient well.

These wells will also provide a representative sampling of the full range of VOC and SVOC concentrations. Table 3-1 shows total concentration of targeted VOCs and SVOCs from the remedial investigation second sampling round. Figures 4-3 and 4-4 depict the distribution of VOCs and SVOCs, respectively, in shallow aquifer zone. These maps and Table 4-1 show that the upgradient and downgradient wells (MW-15S and MW-14S, respectively) feature the lowest organic compound concentration and that the intermediate and high organic compound concentrations are associated with the source area wells (MW-3, MW-6 and MW-11S).

# WEJJEN

# SUMMARY OF VOC AND SVOC ANALYTICAL RESULTS FOR SELECTED GROUNDWATER TREATABILITY STUDY WELLS L.E. CARPENTER SITE, WHARTON, NEW JERSEY

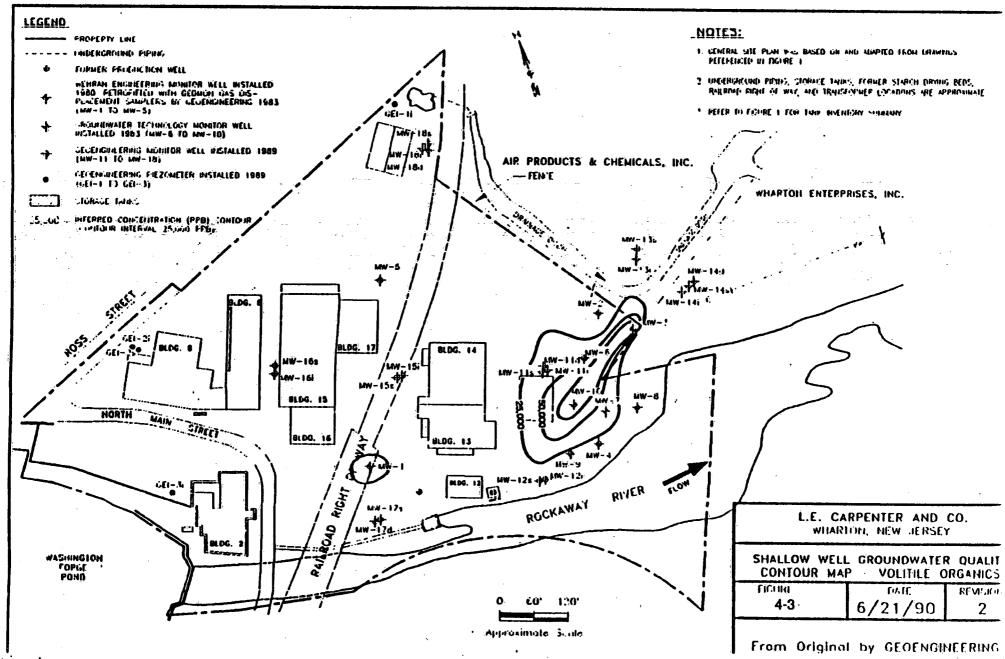
|                      | CONCENTRATION (ug/L) |        |         |        |        |  |  |  |  |  |
|----------------------|----------------------|--------|---------|--------|--------|--|--|--|--|--|
| PARAMETER            | MW-15S               | MW-11S | MW-6    | MW-3   | MW-14S |  |  |  |  |  |
| Total Targeted VOCs  | ND                   | NA     | 136,000 | 39,700 | ND     |  |  |  |  |  |
| Total Targeted SVOCs | ND                   | NA     | 67,898  | 40,723 | 793    |  |  |  |  |  |

<sup>\*</sup> Source: Second groundwater sampling round, WESTON (1990).

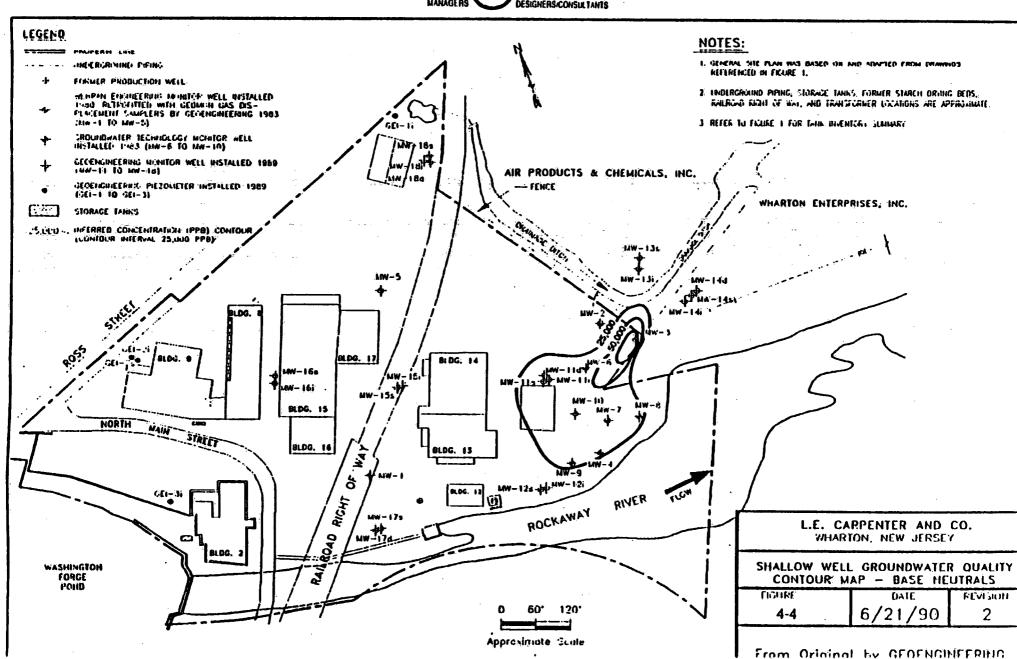
ND = Non Detect

NA = No Analytical Data Available

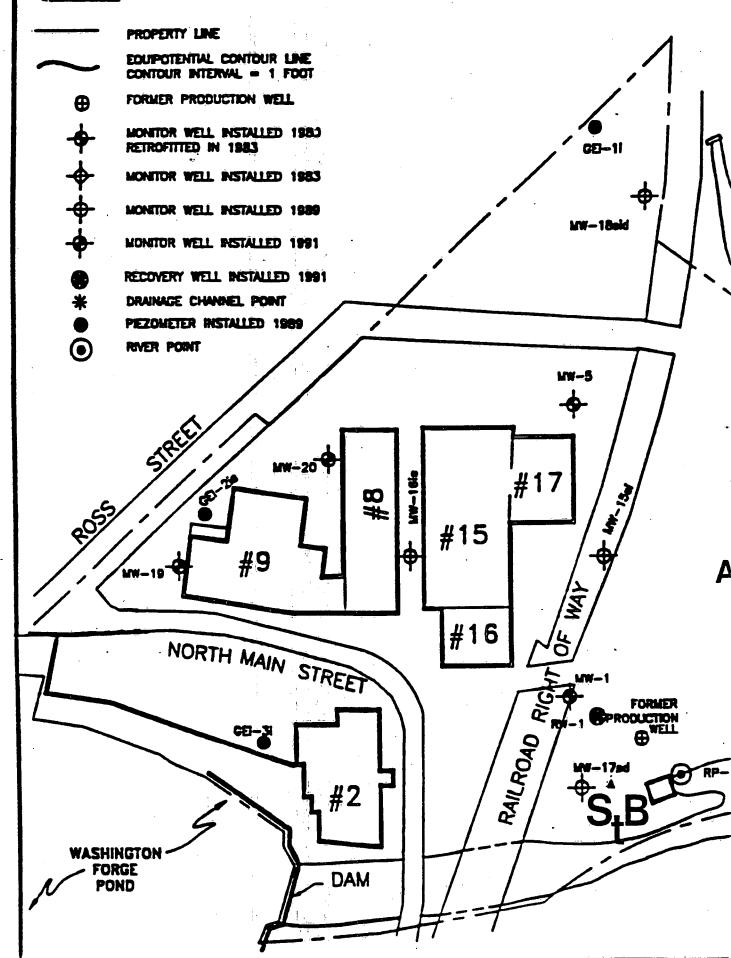


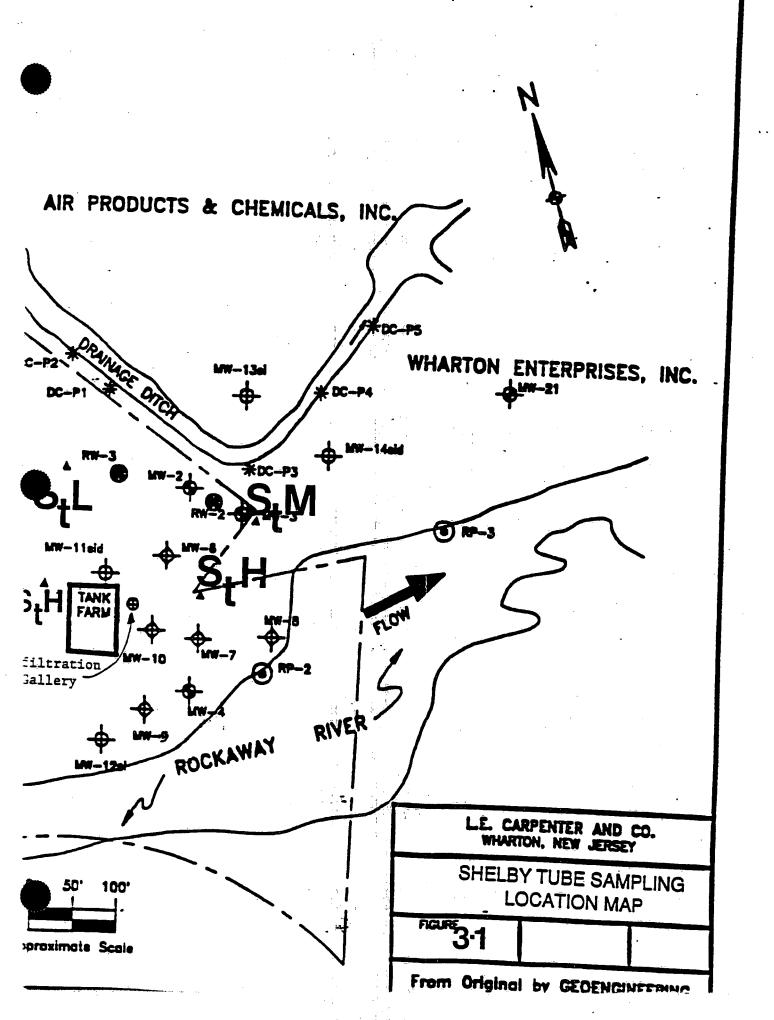




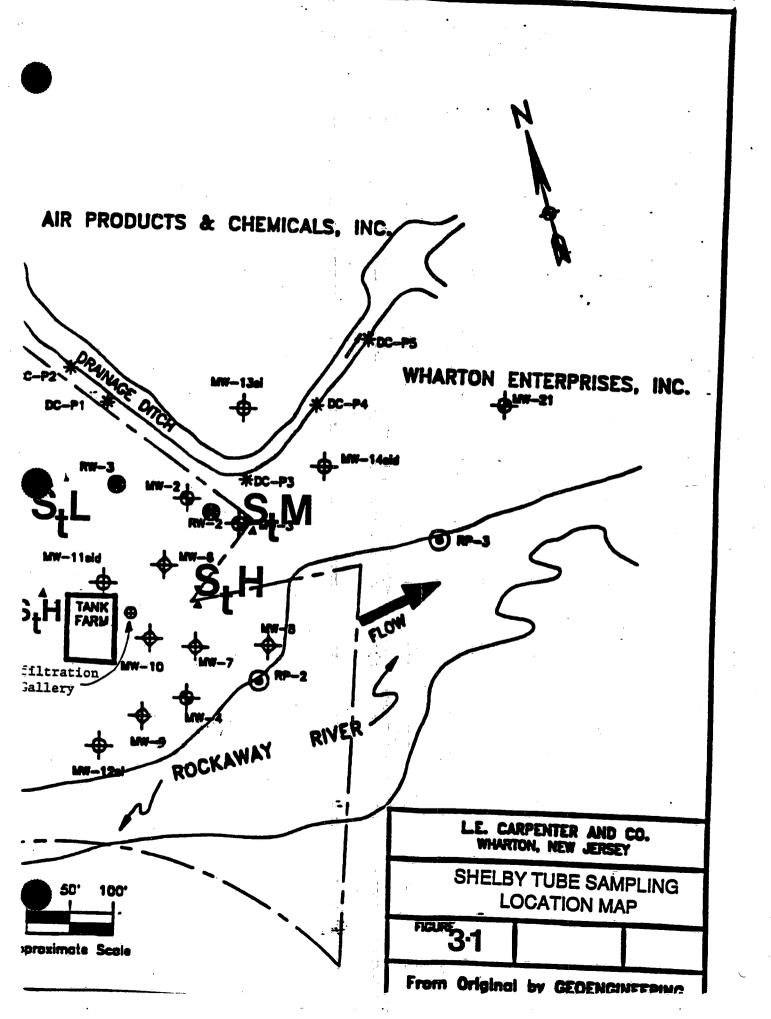


# **LEGEND**





# **LEGEND** PROPERTY LINE EQUIPOTENTIAL CONTOUR LINE CONTOUR DITERVAL - 1 FOOT **(1)** FORMER PRODUCTION WELL MONITOR WELL INSTALLED 1983 RETROFTITED IN 1983 MONITOR WELL INSTALLED 1983 MONITOR WELL INSTALLED 1989 MW-18eld MONITOR WELL INSTALLED 1991 RECOVERY WELL INSTALLED 1991 DRAINAGE CHANNEL POINT PEZOMETER INSTALLED 1989 Ŏ RIVER POINT MW-5 #17 #8 #15 #9 #16 NORTH MAIN STREET Ir RAILROAD RICH FORMER WASHINGTON FORGE DAM POND



# Appendix B

Initial Characterization Data



## ANALYTICAL **SERVICES**

## CERTIFICATE OF ANALYSIS

IT Corporation

312 Directors Drive Knoxville, TN 37923

Attn: Duane Graves

Job Number: REA 50598

March 9, 1992

P.O. Number: 408474.01

This is the Certificate of Analysis for the following samples:

Client Project ID:

Roy F. Weston - L.F. Carpenter

Date Received by Lab:

02/08/92

Number of Samples:

Sample Type:

Six (6) Water

### I. Introduction

On 02/08/92, six (6) water samples arrived at the ITAS-Knoxville, Tennessee laboratory from Roy F. Veston in Edison, New Jersey. The list of analytical tests performed, as well as date of receipt and analysis, can be found in the attached report.

#### II. Analytical Results/Methodology

The analytical results for this report are presented by analytical test. Each set of data will include sample identification information and the analytical results. Please note that the data are not blank corrected.

The samples were analyzed for priority pollutant volatile organic compounds by gas chromatography/mass spectroscopy (GC/MS) based on EPA SW-846 method 8240.

The composite sample was analyzed for the requested metals by inductively coupled plasma spectroscopy (ICP) based on EPA SW-846 method 6010.

Reviewed and Approved:

Robert L. Curry

Laboratory Systems Manager

American Council of Independent Laboratories International Association of Environmental Testing Laboratories American Association for Laboratory Accreditation

Job Number: REA 50598

# II. Analytical Results/Methodology (continued)

Biochemical oxygen demand (BOD) was measured by five day incubation using EPA method 405.1.

Chemical oxygen demand (COD) was measured using the EPA approved HACH procedure, <u>HACH Water Analysis Handbook</u>, HACH Chemical Company, 1980.

The composite sample was analyzed for acidity based on EPA method 305.2.

The composite sample was analyzed for alkalinity based on EPA method 310.3.

The pH of the composite sample was determined using EPA method 150.1.

The composite sample was analyzed for specific conductance based on EPA method 120.1.

The total dissolved solids (TDS) content of the composite sample was determined using EPA method 160.1.

The total solids content of the composite sample was determined using EPA method 160.3.

The total suspended solids (TSS) content of the composite sample was determined using EPA method 160.2.

The composite sample was analyzed for ammonia based on EPA method 350.2.

The composite sample was analyzed for total Kjeldahl nitrogen (TKN) based on EPA method 351.3.

# **III.** Quality Control

Routine laboratory level I QC was followed.

The volatiles analyses were performed by purge and trap with a J&W DB-624 megabore column on a Finnigan OWA GC/MS/DS. The sample runs went well; however, MW3 and MW6 had high levels of ethyl benzene and other non-target organics and exhibited slightly high BFB surrogate recoveries. The dilutions showed normal recovery levels indicating a possible matrix effect from the sample. Ethyl benzene was seen at level below the quantification limit in Method Blank 2 but this had no impact on the validity of the data. There were no problems seen in final data review.

The composite sample was digested on 02/14/92 for ICP and were analyzed by ICP on 02/14/92. All run QC was acceptable. No problems were encountered.

Job Number: REA 50598

#### III. **Quality Control** (continued)

The BOD of the composite sample was measured using a five-day incubation period. The dissolved oxygen (DO) was measured initially and after incubation; the BOD was computed using the difference between the initial and final DO. All run QC was acceptable. No major problems were encountered.

The COD of the composite sample was measured using the reactor digestion method followed by titration with ferrous ammonium sulfate. All run QC was acceptable. No major problems were encountered.

The composite sample was titrated for acidity/alkalinity content. All run QC was acceptable. No major problems were encountered. All run QC was acceptable. No major problems were encountered.

The pH of the composite sample was measured by electrometric procedure; the activity of the hydrogen ions was determined by potentiometric measurement using a standard hydrogen electrode and a reference electrode. All run QC was acceptable. No major problems were encountered.

The composite sample was analyzed for specific conductance. All run QC was acceptable. No major problems were encountered.

The TDS content of the composite sample was determined gravimetrically by filtration and evaporation of the filtrate to a constant weight at 180°C. All run QC was acceptable. No major problems were encountered.

The total solids content of the composite sample was measured gravimetrically by evaporation to dryness at 103-105°C. All run QC was acceptable. No major problems were encountered.

The TSS content of the composite sample was determined gravimetrically by filtration and measurement of the filter residue weight after drying at 103-105°C. All run QC was acceptable. No major problems were encountered.

The composite sample was analyzed for ammonia by distillation followed by Nesslerization using HACH reagents. All run QC was acceptable. No major problems were encountered.

The composite sample was prepared for TKN analysis by acid digestion/distillation using a micro-Kjeldahl system. The distillate was analyzed by Nesslerization using HACH reagents. All run QC was acceptable. No major problems were encountered.

Job Number: REA 50598

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 1 Lab Sample ID: VB02125

| Compound                 | Concentration | Compound                  | Concentration |
|--------------------------|---------------|---------------------------|---------------|
| acrolein                 | 10 U          | 1,1-dichloroethene        | 5 U           |
| acrylonitrile            | 10 U          | trans-1,2-dichloroethene  | 5 U           |
| benzene                  | 5 U           | 1,2-dichloropropane       | 5 U           |
| bromodichloromethane     | 5 U           | cis-1,3-dichloropropene   | 5 U           |
| bromoform                | 5 U           | trans-1,3-dichloropropene | 5 U           |
| bromomethane             | 10 U          | ethyl benzene             | 5 U           |
| carbon tetrachloride     | 5 U           | methylene chloride        | 1 J           |
| chlorobenzene            | 5 U           | 1,1,2,2-tetrachloroethane | 5 U           |
| chloroethane             | 10 U          | tetrachloroethene         | 5 U           |
| 2-chloroethylvinyl ether | 10 U          | toluene                   | 5 Ù           |
| chloroform               | 5 U           | 1,1,1-trichloroethane     | 5 U           |
| chloromethane            | 10 U          | 1,1,2-trichloroethane     | 5 U           |
| dibromochloromethane     | 5 U           | trichloroethene           | 5 U           |
| 1,1-dichloroethane       | <b>5 U</b>    | trichlorofluoromethane    | 5 Ü           |
| 1,2-dichloroethane       | <b>5 U</b>    | vinyl chloride            | 10 U          |

Date of Analysis: 02/12/92

This method blank applies to the following samples: MW2, MW3 and MW6.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Job Number: REA 50598

# **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb)

Sample Matrix: Water

Client Sample ID: MW2 Lab Sample ID: SS5241

| Compound                 | Concent | ration   | Compound                  | Concent | ration |
|--------------------------|---------|----------|---------------------------|---------|--------|
| acrolein                 | 10      | U        | 1,1-dichloroethene        | 5       | U      |
| acrylonitrile            | 10      | U        | trans-1,2-dichloroethene  | .5      | U      |
| benzene                  | .5      | U        | 1,2-dichloropropane       | 5       | U      |
| bromodichloromethane     | 5       | Ü        | cis-1,3-dichloropropene   | .5      | U      |
| bromoform                | .5      | U ·      | trans-1,3-dichloropropene | 5       | U      |
| bromomethane             | 10      | U        | ethyl benzene             | 46      |        |
| carbon tetrachloride     | 5       | U        | methylene chloride        | 5       | U      |
| chlorobenzene            | 5       | U        | 1,1,2,2-tetrachloroethane | 5       | Ų      |
| chloroethane             | 10      | U        | tetrachloroethene         | 5       | U      |
| 2-chloroethylvinyl ether | 10      | U        | toluene                   | 5       | U      |
| chloroform               | 5       | U        | 1,1,1-trichloroethane     | 5       | U      |
| chloromethane            | 10      | <b>U</b> | 1,1,2-trichloroethane     | 5       | U      |
| dibromochloromethane     | 5       | U        | trichloroethene           | 5       | U      |
| 1,1-dichloroethane       | 5       | U        | trichlorofluoromethane    | 5       | U      |
| 1,2-dichloroethane       | 5       | U        | vinyl chloride            | 10      | U      |

Date of Analysis: 02/12/92

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Job Number: REA 50598

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g/liter$  (ppb)

Sample Matrix: Water

Client Sample ID: MW3 Lab Sample ID: SS5242

| Compound                 | Concentration | <u>Compound</u>           | Concent      | ration |
|--------------------------|---------------|---------------------------|--------------|--------|
| acrolein                 | 10 U          | 1,1-dichloroethene        | 5            | U      |
| acrylonitrile            | 10 U          | trans-1,2-dichloroethene  | . 5          | U      |
| benzene                  | 3 Ј           | 1,2-dichloropropane       | 5            | U      |
| bromodichloromethane     | 5 U           | cis-1,3-dichloropropene   | 5            | U      |
| bromoform                | 5 U           | trans-1,3-dichloropropene | 5            | Ù      |
| bromomethane             | 10 U          | ethyl benzene             | 2,000        | D      |
| carbon tetrachloride     | 5 U           | methylene chloride        | 5            | U      |
| chlorobenzene            | 5 U           | 1,1,2,2-tetrachloroethane | 5            | U      |
| chloroethane             | 10 U          | tetrachloroethene         | 5            | Ù      |
| 2-chloroethylvinyl ether | 10 U          | toluene                   | 6            |        |
| chloroform               | 5 U           | 1,1,1-trichloroethane     | <b>. 5</b> . | U      |
| chloromethane            | 10 U          | 1,1,2-trichloroethane     | 5            | U      |
| dibromochloromethane     | 5 U           | trichloroethene           | 5            | U      |
| 1,1-dichloroethane       | 5  U          | trichlorofluoromethane    | 5            | U      |
| 1,2-dichloroethane       | 5 <b>U</b>    | vinyl chloride            | 10           | U      |

Date of Analysis: 02/12/92

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

<sup>-</sup> Indicates an estimated value less than the detection limit.

D - Compound analyzed at a secondary dilution factor.

Job Number: REA 50598

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g/liter$  (ppb)

Sample Matrix: Water

Client Sample ID: MW6 Lab Sample ID: SS5243

| Compound                 | Concentration | Compound                  | Concent | ration |
|--------------------------|---------------|---------------------------|---------|--------|
| acrolein                 | 10 U          | 1,1-dichloroethene        | 5       | U      |
| acrylonitrile            | 10 U          | trans-1,2-dichloroethene  | 5       | Ù      |
| benzene                  | 14            | 1,2-dichloropropane       | 5       | U      |
| bromodichloromethane     | 5 U           | cis-1,3-dichloropropene   | . 5     | U      |
| bromoform                | 5 U           | trans-1,3-dichloropropene | 5       | U      |
| bromomethane             | 10 U          | ethyl benzene             | 13,000  | D      |
| carbon tetrachloride     | 5 U           | methylene chloride        | 5       | U      |
| chlorobenzene            | 5 Ù           | 1,1,2,2-tetrachloroethane | 5       | U      |
| chloroethane             | 10 U          | tetrachloroethene         | 5       | U      |
| 2-chloroethylvinyl ether | 10 U          | toluene                   | 11      |        |
| chloroform               | , 5 U         | 1,1,1-trichloroethane     | 5       | Ü      |
| chloromethane            | 10 U          | 1,1,2-trichloroethane     | 5       | U      |
| dibromochloromethane     | 5 U           | trichloroethene           | 5       | U      |
| 1,1-dichloroethane       | 5 U           | trichlorofluoromethane    | 5       | U      |
| 1,2-dichloroethane       | 5 U           | vinyl chloride            | 3       | J      |

Date of Analysis: 02/12/92

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

<sup>-</sup> Indicates an estimated value less than the detection limit.

D - Compound analyzed at a secondary dilution factor.

KNOXVILLE, TN

Job Number: REA 50598

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in μg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 2 Lab Sample ID: VB02132

| Compound                 | Concentration | Compound                  | Concentration |
|--------------------------|---------------|---------------------------|---------------|
| acrolein                 | 10 U          | 1,1-dichloroethene        | .5 U          |
| acrylonitrile            | 10 U          | trans-1,2-dichloroethene  | 5 U           |
| benzene                  | 5 U           | 1,2-dichloropropane       | . <b>5</b> U  |
| bromodichloromethane     | 5 U           | cis-1,3-dichloropropene   | . <b>5</b> U  |
| bromoform                | 5 U           | trans-1,3-dichloropropene | . <b>5</b> U  |
| bromomethane             | 10 U          | ethyl benzene             | 3 Ј           |
| carbon tetrachloride     | 5 U           | methylene chloride        | .5 U          |
| chlorobenzene            | <b>5</b> U    | 1,1,2,2-tetrachloroethane | 5 U           |
| chloroethane             | 10 U          | tetrachloroethene         | 5 U           |
| 2-chloroethylvinyl ether | 10 U          | toluene                   | <b>5</b> U .  |
| chloroform               | 5 U           | 1,1,1-trichloroethane     | 5 U           |
| chloromethane            | 10 U          | 1,1,2-trichloroethane     | 5 U           |
| dibromochloromethane     | 5 U           | trichloroethene           | 5 U           |
| 1,1-dichloroethane       | 5 <u>U</u>    | trichlorofluoromethane    | 5 U           |
| 1,2-dichloroethane       | 5 U           | vinyl chloride            | 10 U          |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 02/13/92

This method blank applies to the following samples: MW14S, MW15S, MW3 DL and MW6 DL.

DL - Dilution

J - Indicates an estimated value less than the detection limit.

Job Number: REA 50598

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: MW14S Lab Sample ID: SS5244

| Compound                 | Concentration    | Compound                  | Concentration |
|--------------------------|------------------|---------------------------|---------------|
| acrolein                 | 10 U             | 1,1-dichloroethene        | 5 U           |
| acrylonitrile            | 10 U             | trans-1,2-dichloroethene  | , 5 Ü         |
| benzene                  | 5 . U            | 1,2-dichloropropane       | 5 U           |
| bromodichloromethane     | 5 U              | cis-1,3-dichloropropene   | 5 Ü           |
| bromoform                | 5 <sub>\</sub> U | trans-1,3-dichloropropene | 5 U           |
| bromomethane             | 10 U             | ethyl benzene             | 5 U           |
| carbon tetrachloride     | 5 U              | methylene chloride        | 5 U           |
| chlorobenzene            | 5 U              | 1,1,2,2-tetrachloroethane | 5 U           |
| chloroethane             | 10 U             | tetrachloroethene         | 5 U           |
| 2-chloroethylvinyl ether | 10 U             | toluene                   | 5 U           |
| chloroform               | 5 U              | 1,1,1-trichloroethane     | 5 U           |
| chloromethane            | 10 U             | 1,1,2-trichloroethane     | <i>5</i> U    |
| dibromochloromethane     | 5 U              | trichloroethene           | 5 U           |
| 1,1-dichloroethane       | 5 U              | trichlorofluoromethane    | 5 Ü           |
| 1,2-dichloroethane       | 5 U              | vinyl chloride            | 10 U          |

Date of Analysis: 02/13/92

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Job Number: REA 50598

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: MW15S Lab Sample ID: SS5245

| Compound                 | Concentration | Compound                  | Concentration |
|--------------------------|---------------|---------------------------|---------------|
| acrolein                 | 10 U          | 1,1-dichloroethene        | 5 U           |
| acrylonitrile            | 10 U          | trans-1,2-dichloroethene  | 5 U           |
| benzene                  | 5 U           | 1,2-dichloropropane       | 5 U           |
| bromodichloromethane     | 5 U           | cis-1,3-dichloropropene   | 5 Ü           |
| bromoform                | 5 U           | trans-1,3-dichloropropene | 5 U           |
| bromomethane             | 10 U          | ethyl benzene             | 5 U           |
| carbon tetrachloride     | 5 (U          | methylene chloride        | 1. J          |
| chlorobenzene            | 5 U           | 1,1,2,2-tetrachloroethane | 5 U           |
| chloroethane             | 10 U          | tetrachloroethene         | 5 U           |
| 2-chloroethylvinyl ether | 10 U          | toluene                   | 5 U           |
| chloroform               | 5 U           | 1,1,1-trichloroethane     | 5 U           |
| chloromethane            | 10 U          | 1,1,2-trichloroethane     | 5 U           |
| dibromochloromethane     | 5 U           | trichloroethene           | 5 U           |
| 1,1-dichloroethane       | 5 U           | trichlorofluoromethane    | 5 U           |
| 1,2-dichloroethane       | 5 Ù           | vinyl chloride            | 10 U          |

Date of Analysis: 02/13/92

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Job Number: REA 50598

#### WATER SURROGATE PERCENT RECOVERY SUMMARY

|                  |                          | VOLATILE          |                                     |
|------------------|--------------------------|-------------------|-------------------------------------|
| Client Sample ID | Toluene-D8<br>(88-110%)* | BFB<br>(86-115%)* | 1,2 Dichloroethane-D4<br>(76-114%)* |
| MW14S            | 93                       | <b>101</b>        | 94                                  |
| MW15S            | 92                       | 100               | 94                                  |
| MW2              | 95                       | 97                | 92                                  |
| MW3              | 101                      | 125 **            | 95                                  |
| MW3 DL           | 95                       | 103               | 96                                  |
| MW6              | 109                      | 137 **            | 99                                  |
| MW6 DL           | 95                       | 103               | 95                                  |
| Method Blank 1   | 96                       | 96                | 94                                  |
| Method Blank 2   | 95                       | : <b>99</b>       | 94                                  |

DL - Dilution

<sup>\*</sup>Values in parenthesis represent required QC limits.

<sup>\*\*</sup>Values are outside of required QC limits.

Job Number: REA 50598

# **METALS ANALYSIS**

Results in mg/liter (ppm)

Sample Matrix: Water

Client Sample ID: Method Blank Lab Sample ID: PBWE4432

| Compound  | Concentration | Compound  | Concent | ration |
|-----------|---------------|-----------|---------|--------|
| aluminum  | 0.04 U        | lead      | 0.03    | U      |
| antimony  | 0.03 U        | magnesium | 0.03    | U      |
| arsenic   | 0.05 U        | manganese | 0.002   | Ù      |
| barium    | 0.002 U       | nickel    | 0.02    | U      |
| beryllium | 0.001 U       | potassium | 1       | U      |
| cadmium   | 0.005 U       | selenium  | 0.06    | Ü      |
| calcium   | 0.05          | silver    | 0.005   | U      |
| chromium  | 0.01 U        | sodium    | 0.2     | U      |
| cobalt    | 0.02 U        | thallium  | 0.05    | U      |
| copper    | 0.01          | vanadium  | 0.01    | U      |
| iron      | 0.01 U        | zinc      | 0.007   |        |

Date Digested: 02/14/92 Date Analyzed: 02/14/92

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Job Number: REA 50598

#### **METALS ANALYSIS**

Results in mg/liter (ppm)

Sample Matrix: Water

Client Sample ID: COMPOSITE Lab Sample ID: SS5240

| Compound  | Concentration | Compound  | Concentration |
|-----------|---------------|-----------|---------------|
| aluminum  | 15.1          | lead      | 0.03 U        |
| antimony  | 0.03 U        | magnesium | 23.3          |
| arsenic   | 0.05 U        | manganese | 150           |
| barium    | 2.8           | nickel    | 2.4           |
| beryllium | 0.001 U       | potassium | 5             |
| cadmium   | 0.006         | selenium  | 0.09          |
| calcium   | 65.6          | silver    | 0.005 U       |
| chromium  | 3.1           | sodium    | 12.0          |
| cobalt    | 0.15          | thallium  | 0.05 U        |
| copper    | 0.18          | vanadium  | 0.06          |
| iron      | <b>427</b>    | zinc      | 0.23          |

Date Digested: 02/ Date Analyzed: 02/

02/14/92 02/14/92

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Job Number: REA 50598

#### **CLASSICAL PARAMETERS ANALYSIS**

#### Results in mg/liter (ppm) unless otherwise stated

Sample Matrix: Water

| Client Sample ID:                     | Method Blank           |                           |                  |  |
|---------------------------------------|------------------------|---------------------------|------------------|--|
| Lab Sample ID:                        | P3456/P3457/P3476/P350 | 5/P3455/P3464/P3463/P3465 | Date of Analysis |  |
| acidity, as CaCO <sub>3</sub>         | 4                      |                           | 02/10/92         |  |
| alkalinity, as CaCO <sub>3</sub>      | 2                      |                           | 02/10/92         |  |
| ammonia, as N                         | 0.1                    | U                         | 02/13/92         |  |
| biochemical oxygen demand             | +                      |                           | 02/08/92         |  |
| chemical oxygen demand                | •                      |                           | 02/19/92         |  |
| total Kjeldahl nitrogen               | 0.1                    | U                         | 02/20/92         |  |
| pH (standard units)                   | *                      |                           | 02/08/92         |  |
| specific conductance ( $\mu$ mhos/cm) | 4                      |                           | 02/10/92         |  |
| total dissolved solids                | . 8                    |                           | 02/12/92         |  |
| total solids                          | 1                      | U                         | 02/12/92         |  |
| total suspended solids                | 1                      | U                         | 02/12/92         |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

<sup>\* -</sup> A method blank is not applicable for these analyses.

Job Number: REA 50598

# **CLASSICAL PARAMETERS ANALYSIS**

# Results in mg/liter (ppm) unless otherwise stated

Sample Matrix: Water

| Client Sample ID:<br>Lab Sample ID: | COMPOSITE<br>SS5236/SS5237/SS5238/SS5239 | Date of Analysis |
|-------------------------------------|--|------------------|
| acidity, as CaCO <sub>3</sub>       | 86                                       | 02/10/92         |
| alkalinity, as CaCO <sub>3</sub>    | 276                                      | 02/10/92         |
| ammonia, as N                       | 0.7                                      | 02/13/92         |
| biochemical oxygen demand           | 14                                       | 02/08/92         |
| chemical oxygen demand              | <b>57</b>                                | 02/19/92         |
| total Kjeldahl nitrogen             | 3.3                                      | 02/20/92         |
| pH (standard units)                 | 6.30                                     | 02/08/92         |
| specific conductance (µmhos/cm)     | 490                                      | 02/10/92         |
| total dissolved solids              | 290                                      | 02/12/92         |
| total solids                        | 500                                      | 02/12/92         |
| total suspended solids              | 170                                      | 02/12/92         |

# Appendix C

Initial and Final Analytical Data from Batch Groundwater Treatment



# ANALYTICAL SERVICES

# CERTIFICATE OF ANALYSIS

IT Corporation

312 Directors Drive Knoxville, TN 37923 April 7, 1992

Attn: Duane Graves

Job Number: ITEK 50888

P.O. Number: 408474.02

This is the Certificate of Analysis for the following samples:

Client Project ID:

LE Carpenter

Date Received by Lab:

03/11/92

Number of Samples:

Six (6)

Sample Type:

Water

#### I. Introduction

On 03/11/92, six (6) water samples arrived at the ITAS-Knoxville, Tennessee, laboratory from IT 'xporation, Knoxville, Tennessee, in support of the LE Carpenter project. The list of analytical tests performed, as well as date of receipt and analysis, can be found in the attached report.

#### II. Analytical Results/Methodology

The analytical results for this report are presented by analytical test. Each set of data will include sample identification information and the analytical results. Please note that the data are not blank corrected.

The samples were analyzed for priority pollutant volatile organic compounds by gas chromatography/mass spectroscopy (GC/MS) based on EPA SW-846 method 8240.

The samples were analyzed for priority pollutant semivolatile organic compounds by gas chromatography/mass spectroscopy (GC/MS) based on EPA SW-846 method 8270.

Reviewed and Approved:

Alyce R. Moore

Laboratory Manager

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE. TN

Client Project ID: LE Carpenter

Job Number: ITEK 50888

# III. Quality Control

Routine laboratory level I QC was followed.

The volatiles analyses were performed by purge and trap with a J&W DB-624 megabore column on a Finnigan OWA GC/MS/DS. The sample runs went well. Sample TOE and TOD were given additional dilution after encountering some high levels in initial runs, while TOF was run at dilution only because of expected levels. In the TIC analysis, we noted some peaks that were either lab background or obvious carryover in the dilution of sample TOE, and did not report these artifacts. Some analytes reported as TICs were actually quantitated with full standardization. For these compounds, the results were listed without the J qualifier, unless the value was below quantitation limit.

The semivolatiles analyses were performed by direct injection of sample extract on a Restek XTI-5 capillary column on a VG TRIO-2 GC/MS/DS. The sample runs met some problems. Each initial extract run showed a large peak for bis (2-ethylhexyl) phthalate, which appeared to be causing a matrix effect lowering of one internal standard (chrysene-D12) recovery, as well as generally obscuring chromatography in its region. Because of this, manual integration and rechecking for targets were done in that area, and a nearby, unaffected internal standard (phenanthrene-D10) was used as a reference for quantitation. Dilutions were run of the extracts; correcting the matrix effect, and quantitation by the usual internal standard references was possible, and performed. In the TIC analysis, some peaks were "smeared" or split, and the analyst used judgement in distinguishing and omitting peaks belonging to the volatiles fraction. TIC peaks found that had available standardization were reported as for the volatiles. No TICs were found or reported in the dilutions. There were no problems seen in final data review.

Job Number: ITEK 50888

# **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in µg/liter (ppb)
Sample Matrix: Water

Client Sample ID: Method Blank 1 Lab Sample ID: VB03112

| Compound                 | Concent  | ration | Compound                  | Concentration |
|--------------------------|----------|--------|---------------------------|---------------|
| acrolein                 | 10       | U      | 1,1-dichloroethene        | <b>5</b> U    |
| acrylonitrile            | 10       | U      | cis-1,2-dichloroethene    | 5 Ü           |
| benzene                  | 5        | U      | trans-1,2-dichloroethene  | 5 U           |
| bromodichloromethane     | 5        | U      | 1,2-dichloropropane       | 5 U           |
| bromoform                | 5        | U      | cis-1,3-dichloropropene   | .5 U          |
| bromomethane             | 10       | U,     | trans-1,3-dichloropropene | 5 U           |
| carbon tetrachloride     | 5        | U      | ethyl benzene             | 5 U           |
| chlorobenzene            | <b>Ś</b> | U      | methylene chloride        | 3 Ј           |
| chloroethane             | 10       | U      | 1,1,2,2-tetrachloroethane | 5 U           |
| 2-chloroethylvinyl ether | 10       | U      | tetrachloroethene         | 5 U           |
| chloroform               | 5        | U      | toluene                   | 5 U           |
| chloromethane            | 10       | U      | 1,1,1-trichloroethane     | 5 /U          |
| dibromochloromethane     | 5        | U      | 1,1,2-trichloroethane     | 5 U           |
| 1,1-dichloroethane       | . 5      | U      | trichloroethene           | . 5 U         |
| 1,2-dichloroethane       | 5        | U      | trichlorofluoromethane    | 5 U           |
|                          |          | of .   | vinyl chloride            | 10 U          |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 03/11/92

This method blank applies to the following sample: TOD.

J - Indicates an estimated value less than the detection limit.

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

Client Project ID: LE Carpenter

Job Number: ITEK 50888

# ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in  $\mu g$ /liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 1 Lab Sample ID: VB03112

Tentative Identification (1)

Concentration (2)

NONE DETECTED

Remarks:

(1) Identification is based on computer search of the NIST Library.

(2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITEK 50888

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in µg/liter (ppb) Sample Matrix: Water

Client Sample ID: TOD Lab Sample ID: SS7315

| Compound                 | Concentration |   | Compound                  | Concent | ration |
|--------------------------|---------------|---|---------------------------|---------|--------|
| acrolein                 | 10            | U | 1,1-dichloroethene        | .5      | U      |
| acrylonitrile            | 10            | U | cis-1,2-dichloroethene    | 5       | U      |
| benzene                  | 5             |   | trans-1,2-dichloroethene  | 5       | U      |
| bromodichloromethane     | 5             | U | 1,2-dichloropropane       | 5       | U      |
| bromoform                | 5             | U | cis-1,3-dichloropropene   | 5       | U      |
| bromomethane             | 10            | U | trans-1,3-dichloropropene | 5       | U      |
| carbon tetrachloride     | 5             | U | ethyl benzene             | 27,000  | D      |
| chlorobenzene            | <b>5</b> ,    | U | methylene chloride        | 7       | В      |
| chloroethane             | 10            | U | 1,1,2,2-tetrachloroethane | .5      | U      |
| 2-chloroethylvinyl ether | 10            | Ü | tetrachloroethene         | 5       | U      |
| chloroform               | 5             | U | toluene                   | 16      | •      |
| chloromethane            | 10            | Û | 1,1,1-trichloroethane     | . 5     | U      |
| dibromochloromethane     | 5             | U | 1,1,2-trichloroethane     | 5       | U      |
| 1,1-dichloroethane       | 5             | U | trichloroethene           | . 5     | U      |
| 1,2-dichloroethane       | 5             | Ü | trichlorofluoromethane    | .5      | U      |
|                          |               | • | vinyl chloride            | 10      | U      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 03/11/92

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

D - Compound analyzed at a secondary dilution factor.

Job Number: ITEK 50888

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: TOD Lab Sample ID: SS7315

| <u>Tentative Identification (1)</u> |                |     | Concentration (2) |   |  |
|-------------------------------------|----------------|-----|-------------------|---|--|
| 3-pentanone, 2,4-dimethyl           | ļ <del>.</del> |     | 14                |   |  |
| benzene, (1-methylethyl)-           |                |     | 460               |   |  |
| xylenes (total)                     |                | .*  | 32,000            | D |  |
| benzene, propyl-                    |                | * . | 9.0               |   |  |
| benzene, -ethyl-methyl-             |                |     | 73                | Y |  |

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- D Compound analyzed at a secondary dilution factor.
- Y Indistinguishable isomer in tentatively identified compounds.

Job Number: ITEK 50888

# **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: Method Blank 2 Lab Sample ID: VB03122

|   | Compound                 | Concentration |   | Compound                  | Concent | ration |  |
|---|--------------------------|---------------|---|---------------------------|---------|--------|--|
|   | acrolein                 | 10            | U | 1,1-dichloroethene        | 5       | U      |  |
|   | acrylonitrile            | 10            | Ü | cis-1,2-dichloroethene    | 5       | U      |  |
|   | benzene                  | 5             | U | trans-1,2-dichloroethene  | 5       | U      |  |
|   | bromodichloromethane     | 5             | U | 1,2-dichloropropane       | 5       | U      |  |
|   | bromoform                | 5             | U | cis-1,3-dichloropropene   | 5       | U      |  |
|   | bromomethane             | 10            | U | trans-1,3-dichloropropene | 5       | U      |  |
|   | carbon tetrachloride     | <b>5</b> .    | U | ethyl benzene             | 2       | J      |  |
| ŀ | chlorobenzene            | 5             | Ü | methylene chloride        | 2       | J      |  |
|   | chloroethane             | 10            | U | 1,1,2,2-tetrachloroethane | 5       | U      |  |
|   | 2-chloroethylvinyl ether | 10            | U | tëtrachloroëthene         | , 5     | U      |  |
|   | chloroform               | 5             | U | toluene                   | 5       | U      |  |
|   | chloromethane            | 10            | U | 1,1,1-trichloroethane     | 5       | U      |  |
|   | dibromochloromethane     | 5             | U | 1,1,2-trichloroethane     | 5       | Ü      |  |
|   | 1,1-dichloroethane       | 5             | U | trichloroethene           | 5       | U      |  |
|   | 1,2-dichloroethane       | 5             | Ŭ | trichlorofluoromethane    | 5       | U      |  |
|   |                          |               |   | vinyl chloride            | 10      | U      |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 03/12/92

This method blank applies to the following samples: TOD DL and TOE.

DL - Dilution

J - Indicates an estimated value less than the detection limit.

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

Job Number: ITEK 50888

# ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in  $\mu g$ /liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 2 Lab Sample ID: VB03122

Tentative Identification (1)

Concentration (2)

NONE DETECTED

Remarks:

(1) Identification is based on computer search of the NIST Library.

(2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITEK 50888

# **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in µg/liter (ppb)
Sample Matrix: Water

Client Sample ID: TOE Lab Sample ID: SS7316

| Compound                 | <b>Concentration</b> |           | Compound                  | Concent  | ration |
|--------------------------|----------------------|-----------|---------------------------|----------|--------|
| acrolein                 | 10                   | U         | 1,1-dichloroethene        | 5        | U      |
| acrylonitrile            | 10                   | Ŭ         | cis-1,2-dichloroethene    | 5        | U      |
| benzene                  | 6                    |           | trans-1,2-dichloroethene  | 5        | U      |
| bromodichloromethane     | 5                    | U         | 1,2-dichloropropane       | 5        | U      |
| bromoform                | 5                    | T         | cis-1,3-dichloropropene   | 5        | U      |
| bromomethane             | 10                   | U         | trans-1,3-dichloropropene | 5        | U      |
| carbon tetrachloride     | 5                    | <b>'U</b> | ethyl benzene             | 23,000   | DB     |
| chlorobenzene            | 5                    | U         | methylene chloride        | 8        | В      |
| chloroethane             | . 10                 | U         | 1,1,2,2-tetrachloroethane | 5        | U      |
| 2-chloroethylvinyl ether | 10                   | Ü         | tetrachloroethene         | 5        | U      |
| chloroform               | 5                    | U         | toluene                   | 19       |        |
| chloromethane            | 10                   | Ū         | 1,1,1-trichloroethane     | 5        | U      |
| dibromochloromethane     | 5                    | U         | 1,1,2-trichloroethane     | . 5      | U      |
| 1,1-dichloroethane       | 5                    | <b>'U</b> | trichloroethene           | <b>5</b> | U      |
| 1,2-dichloroethane       | 5                    | <b>U</b>  | trichlorofluoromethane    | 5        | U,     |
| ·                        |                      |           | vinyl chloride            | 10       | U      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 03/12/92

J - Indicates an estimated value less than the detection limit.

D - Compound analyzed at a secondary dilution factor.

B - Analyte was found in the blank as well as the sample.

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

'lient Project ID: LE Carpenter

Job Number: ITEK 50888

# ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: TOE Lab Sample ID: SS7316

| Tentative Identification (1) | Concentration (2) |
|------------------------------|-------------------|
| acetone                      | 30                |
| 3-pentanone, 2,4-dimethyl-   | 13                |
| xylenes (total)              | 31,000 D          |
| benzene, methylethyl-        | 230 Y             |

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- D Compound analyzed at a secondary dilution factor.
- Y Indistinguishable isomer in tentatively identified compounds.

Job Number: ITEK 50888

# **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: Method Blank 3

Lab Sample ID: VB0313

| Compound                 | Concentration |                  | Compound                  | Concent | Concentration |  |
|--------------------------|---------------|------------------|---------------------------|---------|---------------|--|
| acrolein                 | 10 T          | ַ ד <u>י</u>     | 1,1-dichloroethene        | .5      | U             |  |
| acrylonitrile            | 10 T          | ת , <sup>י</sup> | cis-1,2-dichloroethene    | 5       | U             |  |
| benzene                  | <b>5</b> T    | U                | trans-1,2-dichloroethene  | 5.      | U             |  |
| bromodichloromethane     | 5 T           | U                | 1,2-dichloropropane       | 5       | U             |  |
| bromoform                | 5 T           | U                | cis-1,3-dichloropropene   | .5      | U             |  |
| bromomethane             | 10 Ţ          | U                | trans-1,3-dichloropropene | 5       | U             |  |
| carbon tetrachloride     | .5 T          | U                | ethyl benzene             | 1       | J ·           |  |
| chlorobenzene            | 5 Ü           | Ü ·              | methylene chloride        | 1       | J             |  |
| chloroethane             | 10 T          | U                | 1,1,2,2-tetrachloroethane | 5       | U             |  |
| 2-chloroethylvinyl ether | 10 T          | Ų                | tetrachloroethene         | 5       | U ·           |  |
| chloroform               | 5 T           | U                | toluene                   | 5       | U             |  |
| chloromethane            | 10 U          | U                | 1,1,1-trichloroethane     | .5      | U             |  |
| dibromochloromethane     | 5 T           | IJ               | 1,1,2-trichloroethane     | 5       | U             |  |
| 1,1-dichloroethane       | 5 T           | U                | trichloroethene           | 5       | U             |  |
| 1,2-dichloroethane       | .5 T          | U                | trichlorofluoromethane    | 5       | Ü             |  |
|                          |               |                  | vinyl chloride            | 10      | U             |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 03/13/92

This method blank applies to the following sample: TOE DL.

DL - Dilution

J - Indicates an estimated value less than the detection limit.

IT ANALYTICAL SERVICES
5815 MIDDLEBROOK PIKE
KNOXVILLE, TN

lient Project ID: LE Carpenter

Job Number: ITEK 50888

# ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 3

Lab Sample ID: VB0313

Tentative Identification (1)

NONE DETECTED

Concentration (2)

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITEK 50888

# **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: Method Blank 4 Lab Sample ID: VB03162

| Compound                 | Concentration |   | Compound                  | Concentrati |     |
|--------------------------|---------------|---|---------------------------|-------------|-----|
| acrolein                 | 10            | U | 1,1-dichloroethene        | 5           | U   |
| acrylonitrile            | 10            | Ų | cis-1,2-dichloroethene    | 5           | U   |
| benzene                  | 5             | U | trans-1,2-dichloroethene  | 5           | U   |
| bromodichloromethane     | 5             | U | 1,2-dichloropropane       | 5           | U   |
| bromoform                | 5             | U | cis-1,3-dichloropropene   | .5          | U   |
| bromomethane             | 10            | U | trans-1,3-dichloropropene | 5           | Ü   |
| carbon tetrachloride     | 5             | U | ethyl benzene             | 2           | J   |
| chlorobenzene            | 5             | U | methylene chloride        | . 1         | j   |
| chloroethane             | 10            | U | 1,1,2,2-tetrachloroethane | .5          | U   |
| 2-chloroethylvinyl ether | 10            | U | tetrachloroethene         | 5           | U . |
| chloroform               | 5             | U | toluene                   | 5           | U   |
| chloromethane            | 10            | U | 1,1,1-trichloroethane     | 5           | U   |
| dibromochloromethane     | 5             | Ū | 1,1,2-trichloroethane     | 5           | U   |
| 1,1-dichloroethane       | 5             | U | trichloroethene           | 5           | U   |
| 1,2-dichloroethane       | 5             | U | trichlorofluoromethane    | 5           | U   |
|                          |               | 4 | vinyl chloride            | 10          | U   |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 03/16/92

This method blank applies to the following sample: TOF.

J - Indicates an estimated value less than the detection limit.

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

Job Number: ITEK 50888

'lient Project ID: LE Carpenter

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 4 Lab Sample ID: VB03162

Tentative Identification (1)

Concentration (2)

xylenes (total)

2

Remarks:

(1) Identification is based on computer search of the NIST Library.

(2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITEK 50888

# **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in µg/liter (ppb) Sample Matrix: Water

Client Sample ID: TOF

Lab Sample ID: SS7317DR4

| Compound                 | Concentratio    | n ¹ Compound              | Concent | ration |
|--------------------------|-----------------|---------------------------|---------|--------|
| acrolein                 | 2,000 U         | 1,1-dichloroethene        | 1,000   | U      |
| acrylonitrile            | 2,000 U         | cis-1,2-dichloroethene    | 1,000   | U      |
| benzene                  | 1, <b>000</b> U | trans-1,2-dichloroethene  | 1,000   | U      |
| bromodichloromethane     | 1,000 U         | 1,2-dichloropropane       | 1,000   | U      |
| bromoform                | 1, <b>000</b> U | cis-1,3-dichloropropene   | 1,000   | U      |
| bromomethane             | 2,000 U         | trans-1,3-dichloropropene | 1,000   | U      |
| carbon tetrachloride     | 1,000 U         | ethyl benzene             | 28,000  | В      |
| chlorobenzene            | 1,000 U         | methylene chloride        | 310     | BJ     |
| chloroethane             | 2,000 U         | 1,1,2,2-tetrachloroethane | 1,000   | U      |
| 2-chloroethylvinyl ether | 2,000 U         | tetrachloroethene         | 1,000   | U ·    |
| chloroform               | 1,000 U         | toluene                   | 1,000   | U      |
| chloromethane            | 2,000 U         | 1,1,1-trichloroethane     | 1,000   | U      |
| dibromochloromethane     | 1,000 U         | 1,1,2-trichloroethane     | 1,000   | U      |
| 1,1-dichloroethane       | 1,000 U         | trichloroethene           | 1,000   | U      |
| 1,2-dichloroethane       | 1,000 U         | trichlorofluoromethane    | 1,000   | U      |
|                          |                 | vinyl chloride            | 2,000   | U      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 03/16/92

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE. TN

Job Number: ITEK 50888

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: TOF

Lab Sample ID: SS7317DR4

Tentative Identification (1)

Concentration (2)

xylenes (total)

42,000 B

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- B Analyte was found in the blank as well as the sample.

Job Number: ITEK 50888

# WATER SURROGATE PERCENT RECOVERY SUMMARY

|                  |                          | VOLATILE          |                                     |
|------------------|--------------------------|-------------------|-------------------------------------|
| Client Sample ID | Toluene-D8<br>(88-110%)* | BFB<br>(86-115%)* | 1,2 Dichloroethane-D4<br>(76-114%)* |
| TOD              | 96                       | 95                | 87                                  |
| TOD DL           | 99                       | 103               | 95                                  |
| TOE              | 101                      | 103               | 89                                  |
| TOE DL           | 89                       | 96                | 98                                  |
| TOF              | 98                       | 103               | 83                                  |
| Method Blank 1   | 91                       | 94                | 89                                  |
| Method Blank 2   | 94                       | 95                | 90                                  |
| Method Blank 3   | 100                      | 97                | 98                                  |
| Method Blank 4   | 96                       | 95                | 84                                  |

DL - Dilution

<sup>\*</sup>Values in parenthesis represent USEPA contract required QC limits.

Job Number: ITEK 50888

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank

Lab Sample ID: H0040

| Compound                   | Concentration | Compound                    | Concentration |
|----------------------------|---------------|-----------------------------|---------------|
| acenaphthene               | 10 U          | bis(2-chloroisopropyl)ether | 10 U          |
| acenaphthylene             | 10 U          | bis(2-ethylhexyl)phthalate  | 10 U          |
| anthracene                 | 10 U          | 4-bromophenyl phenyl ether  | 10 U          |
| benzidine                  | <b>50</b> U   | 2-chloronaphthalene         | 10 U          |
| benzo(a)anthracene         | 10 U          | 4-chlorophenyl phenyl ether | 10 U          |
| benzo(b)fluoranthene       | 10 U          | chrysene                    | 10 U          |
| benzo(k)fluoranthene       | 10 U          | dibenz(a,h)anthracene       | 10 U          |
| benzo(a)pyrene             | <u>10</u> Ü   | di-n-butylphthalate         | 10 Ü          |
| benzo(g,h,i)perylene       | 10 U          | 1,2-dichlorobenzene         | 10 U          |
| butylbenzylphthalate       | 10 U          | 1,3-dichlorobenzene         | 10 U          |
| bis(2-chloroethoxy)methane | 10 U          | 1,4-dichlorobenzene         | 10 U          |
| bis(2-chloroethyl)ether    | 10 U          |                             |               |

Date of Extraction:

03/13/92

Date of Analysis:

03/26/92

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Job Number: ITEK 50888

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank

Lab Sample ID: H0040

| Compound                  | Concentration | Compound                   | Concentration |  |  |
|---------------------------|---------------|----------------------------|---------------|--|--|
| 3,3'dichlorobenzidine     | 20 U          | hexachloroethane           | 10 U          |  |  |
| diethylphthalate          | 10 U          | indeno(1,2,3-cd)pyrene     | 10 U          |  |  |
| dimethyl phthalate        | 10 U          | isophorone                 | 10 U          |  |  |
| 2,4-dinitrotoluene        | 10 U          | naphthalene                | 10 U          |  |  |
| 2,6-dinitrotoluene        | 10 U          | nitrobenzene               | 10 U          |  |  |
| di-n-octyl phthalate      | 10 U          | n-nitroso-di-n-propylamine | 10 U          |  |  |
| 1,2-diphenylhydrazine(1)  | 10 U          | n-nitrosodimethylamine     | 10 U          |  |  |
| fluoranthene              | 10 U          | n-nitrosodiphenylamine(2)  | 10 U          |  |  |
| fluorene                  | 10 U          | phenanthrene               | 10 U          |  |  |
| hexachlorobenzene         | 10 U          | pyrene                     | 10 U          |  |  |
| hexachlorobutadiene       | 10 U          | 1,2,4-trichlorobenzene     | 10 U          |  |  |
| hexachlorocyclopentadiene | 10 U          |                            |               |  |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

- Indicates an estimated value less than the detection limit.

(1) - Screened for as azobenzene (2) - Detected as diphenylamine

Date of Extraction:

03/13/92

Date of Analysis:

03/26/92

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

Job Number: ITEK 50888

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank

Lab Sample ID:

H0040

Tentative Identification (1)

Concentration (2)

unknown

9.8

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITEK 50888

#### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: TOA Lab Sample ID: SS7318

| Compound                   | Concentration | Compound                    | Concentration |
|----------------------------|---------------|-----------------------------|---------------|
| acenaphthene               | 10 U          | bis(2-chloroisopropyl)ether | 10 U          |
| acenaphthylene             | 10 U          | bis(2-ethylhexyl)phthalate  | 54,000 D      |
| anthracene                 | 10 U          | 4-bromophenyl phenyl ether  | 10 U          |
| benzidine                  | 50 U          | 2-chloronaphthalene         | .10 U         |
| benzo(a)anthracene         | 10 U          | 4-chlorophenyl phenyl ether | 10 U          |
| benzo(b)fluoranthene       | 10 U          | chrysene                    | 10 U          |
| benzo(k)fluoranthene       | 10 U          | dibenz(a,h)anthracene       | 10 U          |
| benzo(a)pyrene             | 10 U          | di-n-butylphthalate         | 10 U          |
| benzo(g,h,i)perylene       | 10 U          | 1,2-dichlorobenzene         | 10 U          |
| butylbenzylphthalate       | 4 Ј           | 1,3-dichlorobenzene         | 10 U          |
| bis(2-chloroethoxy)methane | 10 U          | 1,4-dichlorobenzene         | 10 Ü          |
| bis(2-chloroethyl)ether    | 10 U          |                             |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 03/13/92 Date of Analysis:

03/26/92

J Indicates an estimated value less than the detection limit.

Compound analyzed at a secondary dilution factor.

Job Number: ITEK 50888

## BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: TOA Lab Sample ID: SS7318

| Compound                  | Concentration | Compound                   | Concentration |
|---------------------------|---------------|----------------------------|---------------|
| 3,3'dichlorobenzidine     | 20 U          | hexachloroethane           | 10 U          |
| diethylphthalate          | 10 U          | indeno(1,2,3-cd)pyrene     | 10 U          |
| dimethyl phthalate        | 10 U          | isophorone                 | 10 U          |
| 2,4-dinitrotoluene        | 10 U          | naphthalene                | . 10 U        |
| 2,6-dinitrotoluene        | 10 U          | nitrobenzene               | 10 U          |
| di-n-octyl phthalate      | 38            | n-nitroso-di-n-propylamine | 10 Ü          |
| 1,2-diphenylhydrazine(1)  | 10 U          | n-nitrosodimethylamine     | 10 U          |
| fluoranthene              | 10 U          | n-nitrosodiphenylamine(2)  | 10 U          |
| fluorene                  | 10 U          | phenanthrene               | 10 U          |
| hexachlorobenzene         | 10 U          | pyrene                     | 10 U          |
| hexachlorobutadiene       | 10 U          | 1,2,4-trichlorobenzene     | 10 U          |
| hexachlorocyclopentadiene | 10 Ü          |                            |               |

Compound was analyzed for but not detected. The number is the detection limit for the sample.

Indicates an estimated value less than the detection limit.

- Screened for as azobenzene - Detected as diphenylamine

Date of Extraction: Date of Analysis:

03/13/92

03/26/92

Job Number: ITEK 50888

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: TOA Lab Sample ID: SS7318

| Tentative Identification (1) |     | Concentration | <u>n (2)</u> |
|------------------------------|-----|---------------|--------------|
| benzene, (methylethyl)       |     | 140           | Y            |
| benzene, ethyl-methyl-       |     | 47            | Y            |
| benzene, (trimethyl)         |     | 35            | Y            |
| 1-butanol, 3-methyl-, benzoa | 1   | 51            |              |
| ethanone, 1-phenyl-          |     | 32            |              |
| benzene, trimethyl-          | · · | 32            | Y            |
| benzene, trimethyl-          | •   | 27            | Y            |
| propanoic acid, 2-methyl-, 1 |     | 31            |              |
| phosphoric acid, 2-ethylhexy |     | 60            |              |
| 1,2-benzenedicarboxylic acid | •   | 52            |              |
| 1,2-benzenedicarboxylic acid |     | 48            |              |
| 1,2-benzenedicarboxylic acid | •   | 44            |              |
| unknown                      |     | 16            |              |
| 1,2-benzenedicarboxylic acid |     | 42            |              |
| unknown (alkane)             |     | 14            |              |
| 1-decanol, 2-ethyl-          |     | 13            |              |
| unknown                      |     | 13            |              |
| benzenemethanol, .alphaME    |     | 71            |              |
| 2,4-dimethylphenol           |     | 39            |              |
|                              |     |               |              |

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Y - Indistinguishable isomer in tentatively identified compounds.

Job Number: ITEK 50888

#### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: TOB Lab Sample ID: SS7319

| Compound                   | Concent | ration     | Compound                    | Concentr | ation |
|----------------------------|---------|------------|-----------------------------|----------|-------|
| acenaphthene               | 10      | U          | bis(2-chloroisopropyl)ether | 10       | Ų     |
| acenaphthylene             | 10      | U          | bis(2-ethylhexyl)phthalate  | 46,000   | D     |
| anthracene                 | 10      | U          | 4-bromophenyl phenyl ether  | 10       | U     |
| benzidine                  | 50      | U          | 2-chloronaphthalene         | 10       | U     |
| benzo(a)anthracene         | 10      | U          | 4-chlorophenyl phenyl ether | 10       | U     |
| benzo(b)fluoranthene       | 10      | U          | chrysene                    | . 10     | U     |
| benzo(k)fluoranthene       | 10      | U          | dibenz(a,h)anthracene       | 10       | U     |
| benzo(a)pyrene             | 10      | <b>U</b> . | di-n-butylphthalate         | 10       | U     |
| benzo(g,h,i)perylene       | 10      | U          | 1,2-dichlorobenzene         | 10       | U     |
| butylbenzylphthalate       | 10      | U          | 1,3-dichlorobenzene         | 10       | Ü     |
| bis(2-chloroethoxy)methane | 10      | U          | 1,4-dichlorobenzene         | 10       | U     |
| bis(2-chloroethyl)ether    | 10      | U          |                             |          |       |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 03/13/92 Date of Analysis:

03/26/92

Indicates an estimated value less than the detection limit. J

<sup>-</sup> Compound analyzed at a secondary dilution factor.

Job Number: ITEK 50888

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: TOB Lab Sample ID: SS7319

| Compound                  | Concentration | Compound                   | <b>Concentration</b> |
|---------------------------|---------------|----------------------------|----------------------|
| 3,3'dichlorobenzidine     | <b>20</b> U   | hexachloroethane           | 10 Ų                 |
| diethylphthalate          | 10 U          | indeno(1,2,3-cd)pyrene     | . 10 U               |
| dimethyl phthalate        | 10 U          | isophorone                 | 10 U                 |
| 2,4-dinitrotoluene        | 10 U          | naphthalene                | 10 U                 |
| 2,6-dinitrotoluene        | 10 U          | nitrobenzene               | 10 U                 |
| di-n-octyl phthalate      | 17            | n-nitroso-di-n-propylamine | 10 U                 |
| 1,2-diphenylhydrazine(1)  | 10 U          | n-nitrosodimethylamine     | 10 U                 |
| fluoranthene              | 10 Ü          | n-nitrosodiphenylamine(2)  | 10 U                 |
| fluorene                  | 10 <b>U</b>   | phenanthrene               | 10 U                 |
| hexachlorobenzene         | 10 U          | pyrene                     | 10 U                 |
| hexachlorobutadiene       | 10 U          | 1,2,4-trichlorobenzene     | 10 Ú                 |
| hexachlorocyclopentadiene | 10 U          |                            |                      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

(1) - Screened for as azobenzene(2) - Detected as diphenylamine

Date of Extraction:

03/13/92

Date of Analysis:

03/26/92

Job Number: ITEK 50888

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: TOB Lab Sample ID: SS7319

| Tentative Identification (1) | Concentration (2) |
|------------------------------|-------------------|
| benzene, (methylethyl)       | 140 Y             |
| unknown                      | 77                |
| 1,2-benzenedicarboxylic acid | 180               |
| benzene, ethyl-methyl-       | 35 Y              |
| ethanone, 1-phenyl-          | 34                |
| phosphoric acid, 2-ethylhexy | 68                |
| benzene, trimethyl-          | 25 Y              |
| benzene, trimethyl-          | 20 Y              |
| 1-butanol, 3-methyl-, benzoa | 17                |
| propanoic acid, 2-methyl-, 1 | 15                |
| phenol, ethyl-               | 12                |
| 1,2-benzenedicarboxylic acid | 20                |
| 1,2-benzenedicarboxylic acid | 20                |
| 1,2-benzenedicarboxylic acid | 17                |
| 1,2-benzenedicarboxylic acid | 17                |
| unknown (alkane)             | 11                |
| benzenemethanol, .alphaME    | 69                |
| 2,4-dimethylphenol           | 45                |

#### Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Y - Indistinguishable isomer in tentatively identified compounds.

Job Number: ITEK 50888

#### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in  $\mu g$ /liter (ppb)

Sample Matrix: Water

Client Sample ID: TOC Lab Sample ID: SS7320

| Compound                   | Concentra | <u>ion</u> | Compound                    | Concentr | ation |
|----------------------------|-----------|------------|-----------------------------|----------|-------|
| acenaphthene               | 10 U      | J          | bis(2-chloroisopropyl)ether | 10       | Ū     |
| acenaphthylene             | 10 U      | J.         | bis(2-ethylhexyl)phthalate  | 22,000   | Ď     |
| anthracene                 | 10 U      | J          | 4-bromophenyl phenyl ether  | 10       | U     |
| benzidine                  | 50 U      | J          | 2-chloronaphthalene         | 10       | U     |
| benzo(a)anthracene         | 10 U      | J          | 4-chlorophenyl phenyl ether | 10       | U     |
| benzo(b)fluoranthene       | 10 U      | J          | chrysene                    | 10       | U     |
| benzo(k)fluoranthene       | 10 U      | J ·        | dibenz(a,h)anthracene       | 10       | U     |
| benzo(a)pyrene             | 10 · U    | J          | di-n-butylphthalate         | 10       | U     |
| benzo(g,h,i)perylene       | 10 U      | J          | 1,2-dichlorobenzene         | 10       | U     |
| butylbenzylphthalate       | 3 J       |            | 1,3-dichlorobenzene         | . 10     | U     |
| bis(2-chloroethoxy)methane | 10 U      | J.         | 1,4-dichlorobenzene         | 10       | U     |
| bis(2-chloroethyl)ether    | 10 U      | J          |                             |          |       |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 03/13/92 Date of Analysis: 03/26/92

J - Indicates an estimated value less than the detection limit.

D - Compound analyzed at a secondary dilution factor.

Job Number: ITEK 50888

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in  $\mu g/\text{liter (ppb)}$ 

Sample Matrix: Water

Client Sample ID: TOC Lab Sample ID: SS7320

| Compound                  | Concentration | Compound                   | Concentration |
|---------------------------|---------------|----------------------------|---------------|
| 3,3'dichlorobenzidine     | 20 U          | hexachloroethane           | 10 Ų          |
| diethylphthalate          | 10 U          | indeno(1,2,3-cd)pyrene     | 10 U          |
| dimethyl phthalate        | 10 Ü          | isophorone                 | 10 U          |
| 2,4-dinitrotoluene        | 10 Ù          | naphthalene                | 10 U          |
| 2,6-dinitrotoluene        | 10 U          | nitrobenzene               | 10 U          |
| di-n-octyl phthalate      | 30            | n-nitroso-di-n-propylamine | 10 U          |
| 1,2-diphenylhydrazine(1)  | 10 U          | n-nitrosodimethylamine     | 10 U          |
| fluoranthene              | 10 U          | n-nitrosodiphenylamine(2)  | 10 U          |
| fluorene                  | 10 U          | phenanthrene               | 10 U          |
| hexachlorobenzene         | 10 U          | pyrene                     | 10 U          |
| hexachlorobutadiene       | 10 Ü          | 1,2,4-trichlorobenzene     | 10 U          |
| hexachlorocyclopentadiene | 10 U          |                            |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

(1) - Screened for as azobenzene(2) - Detected as diphenylamine

Date of Extraction: 03/13/92 Date of Analysis: 03/26/92

Job Number: ITEK 50888

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: TOC Lab Sample ID: SS7320

| Tentative Identification (1) |            | Concentration | <u>n (2)</u> |
|------------------------------|------------|---------------|--------------|
| benzene, (methylethyl)       |            | 140           | Y            |
| 1,2-benzenedicarboxylic acid |            | 97            |              |
| benzene, ethyl-methyl-       |            | 40            | Y            |
| benzene, trimethyl-          |            | 33            | $\mathbf{Y}$ |
| ethanone, 1-phenyl-          | . t<br>. t | 32            |              |
| benzene, trimethyl-          |            | 31            | Y            |
| 1-butanol, 3-methyl-, benzoa |            | 39            |              |
| propanoic acid, 2-methyl-, 1 |            | 23            |              |
| phosphoric acid, 2-ethylhexy |            | 60            |              |
| benzene, trimethyl-          |            | 23            | Y            |
| 1,2-benzenedicarboxylic acid |            | 52            |              |
| 1,2-benzenedicarboxylic acid |            | 47            |              |
| phenol, ethyl-               |            | 14            | Y            |
| unknown                      |            | 33            |              |
| unknown                      |            | 13            |              |
| unknown (alkane)             |            | 12            |              |
| unknown                      |            | 12            |              |
| phenol, dimethyl-            |            | 9.9           | Y            |
| benzenemethanol, .alphaME    |            | 77            |              |
| 2,4-dimethylphenol           |            | 44            |              |

#### Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Y - Indistinguishable isomer in tentatively identified compounds.

Job Number: ITEK 50888

#### WATER SURROGATE PERCENT RECOVERY SUMMARY

#### **SEMI-VOLATILE**

| Clical Communication | Nitro-<br>Benzene-D5<br>(35-114%)* |       | 2-Fluoro-<br>Biphenyl |              | Terphenyl-<br>D14 |        |                  |
|----------------------|------------------------------------|-------|-----------------------|--------------|-------------------|--------|------------------|
| Client Sample ID     |                                    | (32-1 | 14%)*                 | <u>[43-1</u> | 16%)*             | (33-14 | <del>11%)*</del> |
| TOA                  |                                    | 76    |                       | 90           |                   | 60     |                  |
| TOA DL               |                                    | . 0   | D                     | 0            | D                 | 0      | D                |
| TOB ·                | 1.                                 | 79    |                       | 92           |                   | 54     |                  |
| TOB DL               | v. r                               | : 0   | D                     | 0            | D                 | 0      | D                |
| TOC                  |                                    | 75    |                       | <b>92</b>    | •                 | 54     |                  |
| TOC DL               | . #                                | 0     | D                     | 0            | D                 | 0      | D                |
| Method Blank         |                                    | 75    |                       | 91           |                   | 91     |                  |

<sup>\*</sup>Values in parenthesis represent USEPA contract required QC limits.

D - Surrogates diluted out.



# ANALYTICAL SERVICES

# CERTIFICATE OF ANALYSIS

IT Corporation

312 Directors Drive Knoxville, TN 37923 May 6, 1992

Attn: Duane Graves

Job Number: ITEK 51037

P.O. Number: 408474.02

This is the Certificate of Analysis for the following samples:

Client Project ID:

LE Carpenter

Date Received by Lab:

03/31/92

Number of Samples:

Fifteen (15)

Sample Type:

Water

#### I. Introduction

On 03/31/92, fifteen (15) water samples arrived at the ITAS-Knoxville, Tennessee, laboratory from IT Corporation, Knoxville, Tennessee, in support of the LE Carpenter project. The list of analytical tests performed, as well as date of receipt and analysis, can be found in the attached report.

#### II. Analytical Results/Methodology

The analytical results for this report are presented by analytical test. Each set of data will include sample identification information and the analytical results. Please note that the data are not blank corrected.

The samples were analyzed for priority pollutant volatile organic compounds by gas chromatography/mass spectroscopy (GC/MS) based on EPA SW-846 method 8240.

The samples were analyzed for priority pollutant semivolatile organic compounds by gas chromatography/mass spectroscopy (GC/MS) based on EPA SW-846 method 8270.

Reviewed and Approved:

Alyce R. Moore

aboratory Manager

American Council of Independent Laboratories
International Association of Environmental Testing Laboratories
American Association for Laboratory Accreditation

Job Number: ITEK 51037

#### III. Quality Control

Routine laboratory level I QC was followed.

The volatiles analyses were performed by purge and trap with a J&W DB-624 megabore column on a Finnigan INCOS 500 GC/MS/DS. The sample runs went well. The analysts noted that all bottles received contained headspace. In the TIC analysis, it was found that some analytes (4-methyl-2-pentanone and dimethylbenzene) had been detected, which had been included in the standard. The standard response was used for quantitation in these cases, which produces more accurate values than the usual estimation method. In general, the results were seen to be comparable. The data were reported with J qualifiers as are other TICs. Several samples required additional dilutions. Sample 4A-TF showed a moderate elevation of one surrogate recovery in an undiluted run. This was probably due to matrix effect from high levels. Both of the run data were reported for comparison. Acetonitrile was seen in a method blank, and was recognized as carryover from a standard solvent. This did not interfere with the analysis.

The semivolatiles analyses were performed by direct injection of sample extract on a Restek XTI-5 capillary column on a Finnigan 4500 GC/MS/DS. The sample runs went well. Additional dilutions were necessary to adequately quantitate all analytes. In some cases, terphenyl-D14 surrogate recovery was eported outside acceptable QC limits, but not so as to affect overall sample compliance. In other cases, dilutions required were high enough that surrogates were not reported. For all samples, extraction was begun with 500 ml diluted to 1 liter in reagent water, because of limited volume. The extracts were initially run at twofold bench dilution. Bis-2-ethylhexyl phthalate is a ubiquitous lab contaminant, but the levels seen, and its absence in the method blank, strongly indicate it was sample intrinsic. There were no other problems seen in final data review.

Job Number: ITEK 51037

## **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in µg/liter (ppb)
Sample Matrix: Water

Client Sample ID: 1A-TF Lab Sample ID: SS8682

| Compound                 | Concentration | Compound                  | Concentration |
|--------------------------|---------------|---------------------------|---------------|
| acrolein                 | 10 U          | 1,1-dichloroethene        | 5 U           |
| acrylonitrile            | 10 U          | trans-1,2-dichloroethene  | 5 U           |
| benzene                  | 5. U          | 1,2-dichloropropane       | 5 U           |
| bromodichloromethane     | 5 U           | cis-1,3-dichloropropene   | 5 U           |
| bromoform                | 5 U           | trans-1,3-dichloropropene | 5 U           |
| bromomethane             | 10 Ü          | ethyl benzene             | 5 U           |
| carbon tetrachloride     | 5 U           | methylene chloride        | 4 ЈВ          |
| chlorobenzene            | 5 U           | 1,1,2,2-tetrachloroethane | 5 U           |
| chloroethane             | 10 U          | tetrachloroethene         | 5 U           |
| 2-chloroethylvinyl ether | 10 U          | toluene                   | 5 U           |
| chloroform               | 5 U           | 1,1,1-trichloroethane     | .5 U          |
| chloromethane            | 10 U          | 1,1,2-trichloroethane     | 5 U           |
| dibromochloromethane     | 5 U           | trichloroethene           | 5 U           |
| 1,1-dichloroethane       | 5 U           | trichlorofluoromethane    | 5 U           |
| 1,2-dichloroethane       | 5 U           | vinyl chloride            | 10 U          |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

Client Project ID: LE Carpenter

Job Number: ITEK 51037

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in  $\mu g$ /liter (ppb)

Sample Matrix: Water

Client Sample ID: 1A-TF Lab Sample ID: SS8682

Tentative Identification (1)

Concentration (2)

NONE DETECTED

Remarks:

(1) Identification is based on computer search of the NIST Library.

(2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITEK 51037

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: 1B-TF Lab Sample ID: SS8683R

| Compound                 | Concentrati | ion        | Compound                  | Concent | ration |
|--------------------------|-------------|------------|---------------------------|---------|--------|
| acrolein                 | 10 U        | J          | 1,1-dichloroethene        | 5       | Ü      |
| acrylonitrile            | 10 U        | J          | trans-1,2-dichloroethene  | 5       | U      |
| benzene                  | <b>5</b> U  | J          | 1,2-dichloropropane       | 5       | Û      |
| bromodichloromethane     | 5 U         | J          | cis-1,3-dichloropropene   | 5       | U      |
| bromoform                | 5 U         | j ·        | trans-1,3-dichloropropene | .5      | U      |
| bromomethane             | 10 U        | Ţ          | ethyl benzene             | . 5     | U      |
| carbon tetrachloride     | 5 U         | J          | methylene chloride        | 5       | В      |
| chlorobenzene            | 5 U         | J.         | 1,1,2,2-tetrachloroethane | 5       | U      |
| chloroethane             | 10 U        | J .        | tetrachloroethene         | 5       | U      |
| 2-chloroethylvinyl ether | 10 U        | J          | toluene                   | 5       | U      |
| chloroform               | 5 U         | <b>J</b> : | 1,1,1-trichloroethane     | 5       | U      |
| chloromethane            | 10 U        | J          | 1,1,2-trichloroethane     | 5       | U      |
| dibromochloromethane     | 5 U         | J          | trichloroethene           | 5       | U      |
| 1,1-dichloroethane       | 5 U         | J          | trichlorofluoromethane    | 5       | Ü      |
| 1,2-dichloroethane       | 5 Ų         | J          | vinyl chloride            | 10      | U      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

lient Project ID: LE Carpenter

Job Number: ITEK 51037

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in  $\mu g$ /liter (ppb)

Sample Matrix: Water

Client Sample ID: 1B-TF Lab Sample ID: SS8683

Tentative Identification (1)

NONE DETECTED

Concentration (2)

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITEK 51037

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: 1C-TF Lab Sample ID: SS8684

| Compound                 | Concentration | Compound                  | Concentration |
|--------------------------|---------------|---------------------------|---------------|
| acrolein                 | 10 U          | 1,1-dichloroethene        | 5 U           |
| acrylonitrile            | 10 U          | trans-1,2-dichloroethene  | 5 U           |
| benzene                  | .5 U          | 1,2-dichloropropane       | 5 U           |
| bromodichloromethane     | 5 U           | cis-1,3-dichloropropene   | 5 U           |
| bromoform                | 5 U           | trans-1,3-dichloropropene | 5 U           |
| bromomethane             | 10 Ü          | ethyl benzene             | 5 U           |
| carbon tetrachloride     | .5 U          | methylene chloride        | 5 B           |
| chlorobenzene            | 5 U           | 1,1,2,2-tetrachloroethane | 5 U           |
| chloroethane             | 10 U          | tetrachloroethene         | 5 U           |
| 2-chloroethylvinyl ether | 10 U          | toluene                   | 5 U           |
| chloroform               | 5 U           | 1,1,1-trichloroethane     | 5 U           |
| chloromethane            | 10 U          | 1,1,2-trichloroethane     | 5 U           |
| dibromochloromethane     | 5 U           | trichloroethene           | 5 U           |
| 1,1-dichloroethane       | .5 U          | trichlorofluoromethane    | 5 U           |
| 1,2-dichloroethane       | 5 U           | vinyl chloride            | 10 U          |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

Client Project ID: LE Carpenter

Job Number: ITEK 51037

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1C-TF Lab Sample ID: SS8684

Tentative Identification (1)

Concentration (2)

NONE DETECTED

Remarks:

(1) Identification is based on computer search of the NIST Library.

(2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITEK 51037

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu$ g/liter (ppb) Sample Matrix: Water

Client Sample ID: 2A-TF Lab Sample ID: SS8685

| Compound                 | Concentra | ation | Compound                  | Concent | ration |
|--------------------------|-----------|-------|---------------------------|---------|--------|
| acrolein                 | 10        | U     | 1,1-dichloroethene        | 5       | U      |
| acrylonitrile            | 10        | Ū     | trans-1,2-dichloroethene  | 5       | U      |
| benzene                  | .5        | U     | 1,2-dichloropropane       | 5       | U      |
| bromodichloromethane     | 5         | U     | cis-1,3-dichloropropene   | 5       | U      |
| bromoform                | .5        | U     | trans-1,3-dichloropropene | 5       | U      |
| bromomethane             | 10        | U     | ethyl benzene             | 5       | U      |
| carbon tetrachloride     | 5         | U     | methylene chloride        | 1       | JВ     |
| chlorobenzene            | .5        | U     | 1,1,2,2-tetrachloroethane | 5       | U      |
| chloroethane             | 10        | U,    | tetrachloroethene         | 5       | U      |
| 2-chloroethylvinyl ether | 10        | U     | toluene                   | 5       | U      |
| chloroform               | 1         | J     | 1,1,1-trichloroethane     | 5       | U      |
| chloromethane            | 10        | U     | 1,1,2-trichloroethane     | 5       | U      |
| dibromochloromethane     | 5         | U     | trichloroethene           | 5       | U      |
| 1,1-dichloroethane       | 5         | U     | trichlorofluoromethane    | 5       | U      |
| 1,2-dichloroethane       | .5        | Ų     | vinyl chloride            | 10      | U      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

Client Project ID: LE Carpenter

Job Number: ITEK 51037

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 2A-TF Lab Sample ID: SS8685

Tentative Identification (1)

NONE DETECTED

Concentration (2)

Remarks:

(1) Identification is based on computer search of the NIST Library.

(2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITEK 51037

# **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu$ g/liter (ppb) Sample Matrix: Water

Client Sample ID: 2B-TF Lab Sample ID: SS8686

| Compound                 | Concentration | Compound                  | Concentration |
|--------------------------|---------------|---------------------------|---------------|
| acrolein                 | 10 U          | 1,1-dichloroethene        | 5 U           |
| acrylonitrile            | 10 U          | trans-1,2-dichloroethene  | 5 U           |
| benzene                  | 5 U           | 1,2-dichloropropane       | 5 U           |
| bromodichloromethane     | .5 U          | cis-1,3-dichloropropene   | 5 U           |
| bromoform                | 5   U         | trans-1,3-dichloropropene | .5 U          |
| bromomethane             | 10 U          | ethyl benzene             | 5 U           |
| carbon tetrachloride     | 5 U           | methylene chloride        | 5 U           |
| chlorobenzene            | 5 U           | 1,1,2,2-tetrachloroethane | 5 U           |
| chloroethane             | 10 U          | tetrachloroethene         | .5 U          |
| 2-chloroethylvinyl ether | 10 Ü          | toluëne                   | ,5 U          |
| chloroform               | <b>5 U</b>    | 1,1,1-trichloroethane     | 5 U           |
| chloromethane            | 10 U          | 1,1,2-trichloroethane     | 5 U           |
| dibromochloromethane     | 5 U           | trichloroethene           | ,5 U          |
| 1,1-dichloroethane       | 5 U           | trichlorofluoromethane    | 5 Ų           |
| 1,2-dichloroethane       | 5 U           | vinyl chloride            | 10 U          |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

IT Corporation May 6, 1992

Client Project ID: LE Carpenter

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

Job Number: ITEK 51037

# ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 2B-TF Lab Sample ID: SS8686

Tentative Identification (1)

Concentration (2)

NONE DETECTED

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITEK 51037

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu$ g/liter (ppb) Sample Matrix: Water

Client Sample ID: 2C-TF Lab Sample ID: SS8687

| Compound                 | Concent | ration | Compound                  | Concent  | ration |
|--------------------------|---------|--------|---------------------------|----------|--------|
| acrolein                 | 10      | U      | 1,1-dichloroethene        | 5        | U      |
| acrylonitrile            | 10      | Ü      | trans-1,2-dichloroethene  | 5        | U      |
| benzene                  | ,5      | U      | 1,2-dichloropropane       | 5        | U      |
| bromodichloromethane     | 5       | U      | cis-1,3-dichloropropene   | <b>5</b> | U      |
| bromoform                | 5       | U      | trans-1,3-dichloropropene | 5        | U      |
| bromomethane             | 10      | U      | ethyl benzene             | 5        | U      |
| carbon tetrachloride     | 5       | U      | methylene chloride        | 1        | Љ      |
| chlorobenzene            | 5       | U      | 1,1,2,2-tetrachloroethane | 5        | U      |
| chloroethane             | 10      | U      | tetrachloroethene         | 5        | U      |
| 2-chloroethylvinyl ether | 10      | U      | toluene                   | 5        | U      |
| chloroform               | .5      | U      | 1,1,1-trichloroethane     | 5        | U      |
| chloromethane            | 10      | U      | 1,1,2-trichloroethane     | 5        | U      |
| dibromochloromethane     | 5       | U      | trichloroethene           | 5        | Ü      |
| 1,1-dichloroethane       | 5       | U      | trichlorofluoromethane    | 5        | U      |
| 1,2-dichloroethane       | 5       | Ŭ      | vinyl chloride            | 10       | U      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

Client Project ID: LE Carpenter

Job Number: ITEK 51037

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 2C-TF Lab Sample ID: SS8687

Tentative Identification (1)

Concentration (2)

NONE DETECTED

Remarks:

(1) Identification is based on computer search of the NIST Library.

(2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Client Project ID: LE Carpenter

Job Number: ITEK 51037

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g/liter$  (ppb) Sample Matrix: Water

Client Sample ID: 3A-TF Lab Sample ID: SS8688

| Compound                 | Concentrat | ion | Compound                  | Concent | ratio | Ω |
|--------------------------|------------|-----|---------------------------|---------|-------|---|
| acrolein                 | 10 T       | נו  | 1,1-dichloroethene        | 5       | U     |   |
| acrylonitrile            | 10 T       | Ü   | trans-1,2-dichloroethene  | 5       | Ù     |   |
| benzene                  | 5 τ        | IJ  | 1,2-dichloropropane       | 5       | U     |   |
| bromodichloromethane     | <i>5</i> t | IJ  | cis-1,3-dichloropropene   | 5       | U     |   |
| bromoform                | . 5 T      | ט   | trans-1,3-dichloropropene | . 5     | U     |   |
| bromomethane             | 10 T       | IJ  | ethyl benzene             | 5       | U     |   |
| carbon tetrachloride     | <i>5</i> T |     | methylene chloride        | 3       | JВ    |   |
| chlorobenzene            | 5 T        | Ü   | 1,1,2,2-tetrachloroethane | 5       | U     |   |
| chloroethane             | 10 T       | IJ. | tetrachloroethene         | 5       | U     |   |
| 2-chloroethylvinyl ether | 10 T       | IJ  | toluene                   | 5       | U     |   |
| chloroform               | 5 T        | U   | 1,1,1-trichloroethane     | 5       | U     | ٠ |
| chloromethane            | 10 T       | Ü   | 1,1,2-trichloroethane     | 5       | U     |   |
| dibromochloromethane     | <b>5</b> ( | ט . | trichloroethene           | 5       | U     |   |
| 1,1-dichloroethane       | .5 T       | IJ  | trichlorofluoromethane    | 5       | U     |   |
| 1,2-dichloroethane       | 5 T        | U . | vinyl chloride            | 10      | U     |   |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

Client Project ID: LE Carpenter

Job Number: ITEK 51037

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in  $\mu g/\text{liter}$  (ppb)

Sample Matrix: Water

Client Sample ID: 3A-TF Lab Sample ID: SS8688

Tentative Identification (1)

Concentration (2)

NONE DETECTED

Remarks:

(1) Identification is based on computer search of the NIST Library.

(2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITEK 51037

# VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 3B-TF Lab Sample ID: SS8689

| Compound                 | Concentration | Compound                  | Concent | ration |
|--------------------------|---------------|---------------------------|---------|--------|
| acrolein                 | 10 U          | 1,1-dichloroethene        | 5       | U      |
| acrylonitrile            | 10 U          | trans-1,2-dichloroethene  | 5       | U      |
| benzene                  | 5 U           | 1,2-dichloropropane       | 5       | U      |
| bromodichloromethane     | 5 U           | cis-1,3-dichloropropene   | 5       | U      |
| bromoform                | 5 U           | trans-1,3-dichloropropene | 5       | U      |
| bromomethane             | 10 Ü          | ethyl benzene             | 5       | U      |
| carbon tetrachloride     | 5 U           | methylene chloride        | 4       | JВ     |
| chlorobenzene            | 5 U           | 1,1,2,2-tetrachloroethane | 5       | U      |
| chloroethane             | 10 U          | tetrachloroethene         | 5       | U      |
| 2-chloroethylvinyl ether | 10 U          | toluene                   | 5       | U      |
| chloroform               | 5 U           | 1,1,1-trichloroethane     | 5       | U      |
| chloromethane            | 10 U          | 1,1,2-trichloroethane     | 5       | U      |
| dibromochloromethane     | 5 U           | trichloroethene           | 5       | U      |
| 1,1-dichloroethane       | 5 U           | trichlorofluoromethane    | 5       | U      |
| 1,2-dichloroethane       | 5 U           | vinyl chloride            | 10      | U      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

Client Project ID: LE Carpenter

Job Number: ITEK 51037

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in  $\mu g$ /liter (ppb)

Sample Matrix: Water

Client Sample ID: 3B-TF Lab Sample ID: SS8689

Tentative Identification (1)

Concentration (2)

NONE DETECTED

Remarks:

(1) Identification is based on computer search of the NIST Library.

(2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITEK 51037

# **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: 3C-TF Lab Sample ID: SS8690

| Compound                 | Concentratio      | n <u>Compound</u>         | Concent      | ration   |
|--------------------------|-------------------|---------------------------|--------------|----------|
| acrolein                 | 10 U              | 1,1-dichloroethene        | 5            | Ü        |
| acrylonitrile            | 10 U              | trans-1,2-dichloroethene  | <b>5</b>     | Ü        |
| benzene                  | 5 U               | 1,2-dichloropropane       | , <b>5</b> . | Ū        |
| bromodichloromethane     | <b>5</b>          | cis-1,3-dichloropropene   | 5            | Ų        |
| bromoform                | 5 U               | trans-1,3-dichloropropene | 5            | U        |
| bromomethane             | 10 U              | ethyl benzene             | 5            | U        |
| carbon tetrachloride     | 5 <b>U</b>        | methylene chloride        | 5            | В        |
| chlorobenzene            | 5 <b>U</b>        | 1,1,2,2-tetrachloroethane | 5            | Ü        |
| chloroethane             | 10 <b>U</b>       | tetrachloroethene         | 5            | <b>U</b> |
| 2-chloroethylvinyl ether | 10 U              | toluene                   | 5            | Ü        |
| chloroform               | 5 U               | 1,1,1-trichloroethane     | 5            | U        |
| chloromethane            | 10 U              | 1,1,2-trichloroethane     | 5            | U        |
| dibromochloromethane     | <b>5 U</b>        | trichloroethene           | 5            | Ü        |
| 1,1-dichloroethane       | <i>5</i> <b>U</b> | trichlorofluoromethane    | 5            | U        |
| 1,2-dichloroethane       | 5 U               | vinyl chloride            | 10           | Ü        |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

Client Project ID: LE Carpenter

Job Number: ITEK 51037

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 3C-TF Lab Sample ID: SS8690

Tentative Identification (1)

Concentration (2)

NONE DETECTED

Remarks:

(1) Identification is based on computer search of the NIST Library.

(2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITEK 51037

## **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: 4A-TF Lab Sample ID: \$\$8691

| Compound                 | <u>Concentration</u> | Compound                  | Concent | ration |
|--------------------------|----------------------|---------------------------|---------|--------|
| acrolein                 | 10 U                 | 1,1-dichloroethene        | 5       | U      |
| acrylonitrile            | 10 U                 | trans-1,2-dichloroethene  | 5       | U      |
| benzene                  | 400 E                | 1,2-dichloropropane       | 5       | U      |
| bromodichloromethane     | 5 U                  | cis-1,3-dichloropropene   | 5       | TU.    |
| bromoform                | 5 <b>U</b>           | trans-1,3-dichloropropene | 5       | U      |
| bromomethane             | 10 Ù                 | ethyl benzene             | 6,500   | D      |
| carbon tetrachloride     | 5 U                  | methylene chloride        | 7       | B      |
| chlorobenzene            | 160                  | 1,1,2,2-tetrachloroethane | 5       | U      |
| chloroethane             | 10 U                 | tetrachloroethene         | . 5     | Ú      |
| 2-chloroethylvinyl ether | 10 U                 | toluene                   | 8       |        |
| chloroform               | 5 U                  | 1,1,1-trichloroethane     | ·. 5    | U      |
| chloromethane            | 10 U                 | 1,1,2-trichloroethane     | . 5     | Ų      |
| dibromochloromethane     | <b>5</b> Ü           | trichloroethene           | 5       | U      |
| 1,1-dichloroethane       | 5 U                  | trichlorofluoromethane    | 5       | U      |
| 1,2-dichloroethane       | 5 U                  | vinyl chloride            | 10      | U.     |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

D - Compound analyzed at a secondary dilution factor.

E - Compound exceeded calibration range of instrument.

lient Project ID: LE Carpenter

Job Number: ITEK 51037

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 4A-TF Lab Sample ID: SS8691

| Tentative Identification (1)     |  | Concentration (2) |    |  |
|----------------------------------|--|-------------------|----|--|
| 4-methyl-2-pentanone             |  | 400               | J. |  |
| benzene, dimethyl                |  | 1,000             | JΥ |  |
| benzene, 1,2-dimethyl            |  | 4,100             | 1  |  |
| 1,3,5-cycloheptatriene, 7-ethyl- |  | 120               | J  |  |
| 1,3,5-cycloheptatriene, 7-ethyl- |  | 16                | J  |  |

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- J Indicates an estimated value less than the detection limit.
- Y Indistinguishable isomer in tentatively identified compounds.

Job Number: ITEK 51037

# **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: 4B-TF Lab Sample ID: SS8692

| Compound                 | Concentration     | Compound                  | Concentration |
|--------------------------|-------------------|---------------------------|---------------|
| acrolein                 | 10 <sup>±</sup> U | 1,1-dichloroethene        | .5 U          |
| acrylonitrile            | 10 U              | trans-1,2-dichloroethene  | <b>5</b> Ü    |
| benzene                  | 54                | 1,2-dichloropropane       | 5 U           |
| bromodichloromethane     | 5 U               | cis-1,3-dichloropropene   | 5 U           |
| bromoform                | 5 U               | trans-1,3-dichloropropene | 5 U           |
| bromomethane             | 10 U              | ethyl benzene             | 3,600 D       |
| carbon tetrachloride     | 5 U               | methylene chloride        | 7             |
| chlorobenzene            | 70                | 1,1,2,2-tetrachloroethane | 5 Ü           |
| chloroethane             | 10 U              | tetrachloroethene         | 5 U           |
| 2-chloroethylvinyl ether | 10 U              | toluëne                   | 5             |
| chloroform               | 5 U               | 1,1,1-trichloroethane     | 5 U           |
| chloromethane            | 10 U              | 1,1,2-trichloroethane     | 5 U           |
| dibromochloromethane     | <b>5</b> U        | trichloroethene           | 5 U           |
| 1,1-dichloroethane       | <i>5</i> U        | trichlorofluoromethane    | 5 U           |
| 1,2-dichloroethane       | 5 U               | vinyl chloride            | 10 U          |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

<sup>-</sup> Indicates an estimated value less than the detection limit.

D - Compound analyzed at a secondary dilution factor.

Client Project ID: LE Carpenter

Job Number: ITEK 51037

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 4B-TF Lab Sample ID: SS8692

| Tentative Identification (1)     | <u>Concentra</u> | tion (2) |
|----------------------------------|------------------|----------|
| 3-pentanone, 2,4-dimethyl        | 4                | J        |
| benzene, dimethyl                | 750              | JY       |
| benzene, 1,2-dimethyl            | 2,800            | J        |
| 1,3,5-cycloheptatriene, 7-ethyl- | 24               | J        |
| 1,3,5-cycloheptatriene, 7-ethyl- | 3                | J        |

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- J Indicates an estimated value less than the detection limit.
- Y Indistinguishable isomer in tentatively identified compounds.

Job Number: ITEK 51037

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in µg/liter (ppb) Sample Matrix: Water

Client Sample ID: 4C-TF Lab Sample ID: SS8693

| Compound                 | <b>Concentration</b> | Compound                  | Concent | ration |
|--------------------------|----------------------|---------------------------|---------|--------|
| acrolein                 | 100 U                | 1,1-dichloroethene        | 50      | U      |
| acrylonitrile            | 100 U                | trans-1,2-dichloroethene  | 50      | U      |
| benzene                  | 250                  | 1,2-dichloropropane       | 50      | U      |
| bromodichloromethane     | 50 U                 | cis-1,3-dichloropropene   | 50      | U      |
| bromoform                | <i>5</i> 0 U         | trans-1,3-dichloropropene | 50      | U      |
| bromomethane             | 100 U                | ethyl benzene             | 7,500   | D      |
| carbon tetrachloride     | 50 Ü                 | methylene chloride        | 16      | JВ     |
| chlorobenzene            | 96 J                 | 1,1,2,2-tetrachloroethane | 50      | U      |
| chloroethane             | 100 U                | tetrachloroethene         | 50      | U      |
| 2-chloroethylvinyl ether | 100 U                | toluene                   | 50      | U      |
| chloroform               | 50 U                 | 1,1,1-trichloroethane     | 50      | Ų      |
| chloromethane            | 1 <b>00</b> U        | 1,1,2-trichloroethane     | 50      | U      |
| dibromochloromethane     | <b>50 U</b>          | trichloroethene           | 50      | U      |
| 1,1-dichloroethane       | 50 U                 | trichlorofluoromethane    | 50      | U      |
| 1,2-dichloroethane       | <b>50</b> U          | vinyl chloride            | 100     | Ü.     |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

D - Compound analyzed at a secondary dilution factor.

Client Project ID: LE Carpenter

Job Number: ITEK 51037

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 4C-TF Lab Sample ID: SS8693

Tentative Identification (1)

Concentration (2)

benzene, dimethyl

690 JY

benzene, 1,2-dimethyl

8,700 J

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- J Indicates an estimated value less than the detection limit.
- Y Indistinguishable isomer in tentatively identified compounds.

Job Number: ITEK 51037

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g/liter$  (ppb) Sample Matrix: Water

Client Sample ID: 5A-TF Lab Sample ID: SS8694

| Compound                 | Concentra | ation | Compound                  | Concer | tration |
|--------------------------|-----------|-------|---------------------------|--------|---------|
| acrolein                 | 100       | Ü     | 1,1-dichloroethene        | 50     | U       |
| acrylonitrile            | 100       | U     | trans-1,2-dichloroethene  | 50     | U       |
| benzene                  | 50        | U     | 1,2-dichloropropane       | 50     | U       |
| bromodichloromethane     | 50        | U     | cis-1,3-dichloropropene   | 50     | U       |
| bromoform                | 50        | U     | trans-1,3-dichloropropene | 50     | U       |
| bromomethane             | 100       | U     | ethyl benzene             | 14,000 | D       |
| carbon tetrachloride     | 50        | U     | methylene chloride        | 18     | JB      |
| chlorobenzene            | 50        | U     | 1,1,2,2-tetrachloroethane | 50     | U       |
| chloroethane             | 100       | Ü     | tetrachloroethene         | 50     | U       |
| 2-chloroethylvinyl ether | 100       | U     | toluene                   | 11     | J       |
| chloroform               | 50        | U     | 1,1,1-trichloroethane     | 50     | U       |
| chloromethane            | 100       | U     | 1,1,2-trichloroethane     | 50     | U       |
| dibromochloromethane     | 50        | U     | trichloroethene           | 50     | U       |
| 1,1-dichloroethane       | 50        | Ü     | trichlorofluoromethane    | 50     | U       |
| 1,2-dichloroethane       | 50        | U     | vinyl chloride            | 100    | U       |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

D - Compound analyzed at a secondary dilution factor.

Client Project ID: LE Carpenter

Job Number: ITEK 51037

# ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 5A-TF Lab Sample ID: SS8694

Tentative Identification (1)

benzene, dimethyl

benzene, 1,2-dimethyl

1,800 JY

14,000 J

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- J Indicates an estimated value less than the detection limit.
- Y Indistinguishable isomer in tentatively identified compounds.

Job Number: ITEK 51037

# **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in µg/liter (ppb) Sample Matrix: Water

Client Sample ID: 5B-TF Lab Sample ID: SS8695

| Compound                 | Concent | ration | Compound                  | Concent | ration     |
|--------------------------|---------|--------|---------------------------|---------|------------|
| acrolein                 | 100     | Ù      | 1,1-dichloroethene        | 50      | U          |
| acrylonitrile            | 100     | U      | trans-1,2-dichloroethene  | 50      | U          |
| benzene                  | 50      | U      | 1,2-dichloropropane       | 50      | U          |
| bromodichloromethane     | 50      | U      | cis-1,3-dichloropropene   | 50      | U          |
| bromoform                | 50      | U      | trans-1,3-dichloropropene | 50      | U          |
| bromomethane             | 100     | Ŭ      | ethyl benzene             | 15,000  | D          |
| carbon tetrachloride     | 50      | U      | methylene chloride        | 19      | ЛВ         |
| chlorobenzene            | 50      | U      | 1,1,2,2-tetrachloroethane | 50      | U          |
| chloroethane             | 100     | U,     | tetrachloroethene         | 50      | U          |
| 2-chloroethylvinyl ether | 100     | U      | toluene                   | 13      | <b>J</b> · |
| chloroform               | 50      | U      | 1,1,1-trichloroethane     | 50      | U          |
| chloromethane            | 100     | U      | 1,1,2-trichloroethane     | 50      | U          |
| dibromochloromethane     | 50      | U      | trichloroethene           | 50      | U          |
| 1,1-dichloroethane       | 50      | U      | trichlorofluoromethane    | . 50    | U          |
| 1,2-dichloroethane       | 50      | U      | vinyl chloride            | 100     | U          |

Date of Analysis: 04/03/92

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

D - Compound analyzed at a secondary dilution factor.

Client Project ID: LE Carpenter

Job Number: ITEK 51037

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 5B-TF Lab Sample ID: SS8695

Tentative Identification (1)

Concentration (2)

benzene, dimethyl

1,900 JY

benzene, 1,2-dimethyl

14,000 J

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- J Indicates an estimated value less than the detection limit.
- Y Indistinguishable isomer in tentatively identified compounds.

Job Number: ITEK 51037

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: 5C-TF (1:100 Dilution)

Lab Sample ID: SS8696

| Compound                 | Concentration   | Compound                  | Concentration |
|--------------------------|-----------------|---------------------------|---------------|
| acrolein                 | 1,000 U         | 1,1-dichloroethene        | 500 U         |
| acrylonitrile            | 1,000 U         | trans-1,2-dichloroethene  | <b>500</b> U  |
| benzene                  | 500 JU          | 1,2-dichloropropane       | <b>500</b> U  |
| bromodichloromethane     | 500 U           | cis-1,3-dichloropropene   | 500 U         |
| bromoform                | 500 U           | trans-1,3-dichloropropene | <b>500</b> U  |
| bromomethane             | 1,000 U         | ethyl benzene             | 15,000        |
| carbon tetrachloride     | 500 U           | methylene chloride        | 210 J         |
| chlorobenzene            | 500 U           | 1,1,2,2-tetrachloroethane | <b>500</b> U  |
| chloroethane             | 1,000 U         | tetrachloroethene         | <b>500</b> U  |
| 2-chloroethylvinyl ether | 1,000 U         | toluene                   | <b>500</b> U  |
| chloroform               | 500 U           | 1,1,1-trichloroethane     | 500 U         |
| chloromethane            | 1, <b>000</b> U | 1,1,2-trichloroethane     | <b>500</b> U  |
| dibromochloromethane     | 500 Ŭ           | trichloroethene           | <b>500</b> U  |
| 1,1-dichloroethane       | .500 U          | trichlorofluoromethane    | 500 U         |
| 1,2-dichloroethane       | 500 U           | vinyl chloride            | 1,000 U       |

Date of Analysis: 04/06/92

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Client Project ID: LE Carpenter

Job Number: ITEK 51037

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 5C-TF (1:100 Dilution)

Lab Sample ID: SS8696

Tentative Identification (1)

benzene, dimethyl
benzene, 1,2-dimethyl
25,000 Y

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- Y Indistinguishable isomer in tentatively identified compounds.

Job Number: ITEK 51037

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: Method Blank 1

Lab Sample ID: WB0402

| Co  | mpound                | Concent | ration | Compound                  | Concent | ration |
|-----|-----------------------|---------|--------|---------------------------|---------|--------|
| acr | olein                 | 10      | U      | 1,1-dichloroethene        | 5       | U      |
| acr | ylonitrile            | 10      | U      | trans-1,2-dichloroethene  | 5       | U      |
| ber | ızene                 | 5       | U      | 1,2-dichloropropane       | 5       | U      |
| bro | modichloromethane     | 5       | U      | cis-1,3-dichloropropene   | 5       | U      |
| bro | omoform               | 5       | U      | trans-1,3-dichloropropene | . 5     | U      |
| bro | omomethane            | 10      | U      | ethyl benzene             | 5       | Ü      |
| car | bon tetrachloride     | 5       | U      | methylene chloride        | 1       | J      |
| chl | orobenzene            | 5       | U      | 1,1,2,2-tetrachloroethane | 5       | U      |
| chi | oroethane             | 10      | U      | tetrachloroethene         | 5       | U      |
| 2-c | hloroethylvinyl ether | 10      | U      | toluene                   | 5       | U      |
| chi | oroform               | 5       | U.     | 1,1,1-trichloroethane     | 5       | U      |
| chi | oromethane            | 10      | U      | 1,1,2-trichloroethane     | 5       | U      |
| dib | romochloromethane     | 5       | U      | trichloroethene           | 5       | U      |
| 1,1 | -dichloroethane       | 5       | U      | trichlorofluoromethane    | 5       | U      |
| 1,2 | -dichloroethane       | :5      | U      | vinyl chloride            | 10      | U      |

Date of Analysis: 04/02/92

This method blank applies to the following samples: 1A-TF, 1B-TF, 1C-TF, 2A-TF, 2B-TF, 2C-TF and 3A-TF.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Client Project ID: LE Carpenter

Job Number: ITEK 51037

# ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 1

Lab Sample ID: WB0402

Tentative Identification (1)

Concentration (2)

NONE DETECTED

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITEK 51037

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in µg/liter (ppb)
Sample Matrix: Water

Client Sample ID: Method Blank 2 Lab Sample ID: WB04023

| Compound                 | Concentration | Compound                  | Concentration |
|--------------------------|---------------|---------------------------|---------------|
| acrolein                 | 10 U          | 1,1-dichloroethene        | 5 U           |
| acrylonitrile            | 10 U          | trans-1,2-dichloroethene  | 5 U           |
| benzene                  | 5 U           | 1,2-dichloropropane       | 5 U           |
| bromodichloromethane     | 5 ុប          | cis-1,3-dichloropropene   | 5 U           |
| bromoform                | 5 U           | trans-1,3-dichloropropene | 5 U           |
| bromomethane             | 10 U          | ethyl benzene             | 5 U           |
| carbon tetrachloride     | 5 U           | methylene chloride        | 1 J           |
| chlorobenzene            | 5 U           | 1,1,2,2-tetrachloroethane | 5 U           |
| chloroethane             | 10 U          | tetrachloroethene         | .5 U          |
| 2-chloroethylvinyl ether | 10 U          | toluene                   | 5 U           |
| chloroform               | 5 U           | 1,1,1-trichloroethane     | 5 U           |
| chloromethane            | 10 Ü          | 1,1,2-trichloroethane     | 5 U           |
| dibromochloromethane     | 5 U           | trichloroethene           | 5 U           |
| 1,1-dichloroethane       | 5 U           | trichlorofluoromethane    | :5 U          |
| 1,2-dichloroethane       | 5 U           | vinyl chloride            | 10 U          |

Date of Analysis: 04/02/92

This method blank applies to the following samples: 3B-TF, 3C-TF and 4A-TF.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Client Project ID: LE Carpenter

Job Number: ITEK 51037

# ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 2 Lab Sample ID: WB04023

Tentative Identification (1)

Concentration (2)

NONE DETECTED

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITEK 51037

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g/liter$  (ppb) Sample Matrix: Water

Client Sample ID: Method Blank 3

Lab Sample ID: WB0403

| Compound                 | Concentration       | Compound                  | Concentration |
|--------------------------|---------------------|---------------------------|---------------|
| acrolein                 | 10 U                | 1,1-dichloroethene        | 5 U           |
| acrylonitrile            | 10 U                | trans-1,2-dichloroethene  | 5 U           |
| benzene                  | <b>5</b> • <b>U</b> | 1,2-dichloropropane       | 5 U           |
| bromodichloromethane     | 5 U                 | cis-1,3-dichloropropene   | 5 U           |
| bromoform                | 5 U                 | trans-1,3-dichloropropene | 5 U           |
| bromomethane             | 10 U                | ethyl benzene             | 5 U           |
| carbon tetrachloride     | 5 U                 | methylene chloride        | 5 U           |
| chlorobenzene            | 5 U                 | 1,1,2,2-tetrachloroethane | 5 U           |
| chloroethane             | 10 U                | tetrachloroethene         | , <b>5 U</b>  |
| 2-chloroethylvinyl ether | 10 U                | toluene                   | 5 U           |
| chloroform               | 5 'U                | 1,1,1-trichloroethane     | 5 U           |
| chloromethane            | 10 U                | 1,1,2-trichloroethane     | 5 U           |
| dibromochloromethane     | 5 U                 | trichloroethene           | 5 U           |
| 1,1-dichloroethane       | 5 U                 | trichlorofluoromethane    | 5 U           |
| 1,2-dichloroethane       | 5 U                 | vinyl chloride            | 10 U          |

Date of Analysis: 04/03/92

This method blank applies to the following samples: 4A-TF DL, 4A-TF DL2, 4B-TF and 4C-TF DL.

DL = Dilution

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Client Project ID: LE Carpenter

Job Number: ITEK 51037

# ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 3

Lab Sample ID: WB0403

Tentative Identification (1)

Concentration (2)

NONE DETECTED

Remarks:

(1) Identification is based on computer search of the NIST Library.

(2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITEK 51037

# **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g/liter$  (ppb) Sample Matrix: Water

Client Sample ID: Method Blank 4 Lab Sample ID: WB04032

| Compound                 | Concentration | on Compound               | Concent | ration |
|--------------------------|---------------|---------------------------|---------|--------|
| acrolein                 | 10 U          | 1,1-dichloroethene        | 5       | U      |
| acrylonitrile            | 10 U          | trans-1,2-dichloroethene  | 5       | Ü      |
| benzene                  | 5 U           | 1,2-dichloropropane       | 5       | U      |
| bromodichloromethane     | 5 U           | cis-1,3-dichloropropene   | 5       | U      |
| bromoform                | 5 U           | trans-1,3-dichloropropene | 5       | Ū      |
| bromomethane             | 10 U          | ethyl benzene             | 5       | U      |
| carbon tetrachloride     | <b>5 U</b> .  | methylene chloride        | 1       | J      |
| chlorobenzene            | 5 U           | 1,1,2,2-tetrachloroethane | 5       | U      |
| chloroethane             | 10 .U         | tetrachloroethene         | 5       | U      |
| 2-chloroethylvinyl ether | 10 U          | toluene                   | 5       | U      |
| chloroform               | 5 U           | 1,1,1-trichloroethane     | 5       | U      |
| chloromethane            | 10 U          | 1,1,2-trichloroethane     | 5       | U      |
| dibromochloromethane     | 5 U           | trichloroethene           | 5       | U      |
| 1,1-dichloroethane       | 5 U           | trichlorofluoromethane    | 5       | U      |
| 1,2-dichloroethane       | 5 U           | vinyl chloride            | 10      | Ü      |

Date of Analysis: 04/03/92

This method blank applies to the following samples: 4B-TF DL, 4C-TF, 5A-TF, 5A-TF DL, 5B-TF and 5B-TF DL.

DL = Dilution

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

lient Project ID: LE Carpenter

Job Number: ITEK 51037

# ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in  $\mu g/liter$  (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 4 Lab Sample ID: WB04032

Tentative Identification (1)

Concentration (2)

NONE DETECTED

Remarks:

(1) Identification is based on computer search of the NIST Library.

(2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITEK 51037

# **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: Method Blank 5

Lab Sample ID: WB0406

| Compound                 | Concentration       | Compound                  | Concent | ration |
|--------------------------|---------------------|---------------------------|---------|--------|
| acrolein                 | 10 U                | 1,1-dichloroethene        | 5       | U      |
| acrylonitrile            | 10 U                | trans-1,2-dichloroethene  | 5       | U      |
| benzene                  | 5 U                 | 1,2-dichloropropane       | 5       | U      |
| bromodichloromethane     | 5 U                 | cis-1,3-dichloropropene   | 5       | U      |
| bromoform                | .5 U                | trans-1,3-dichloropropene | 5       | U      |
| bromomethane             | 10 U                | ethyl benzene             | . 5     | U      |
| carbon tetrachloride     | 5 U                 | methylene chloride        | 5       | Ų      |
| chlorobenzene            | 5 U                 | 1,1,2,2-tetrachloroethane | 5       | U      |
| chloroethane             | 10 U                | tetrachloroethene         | 5       | Ü      |
| 2-chloroethylvinyl ether | 10 U                | toluene                   | 5       | U      |
| chloroform               | 5 Ü                 | 1,1,1-trichloroethane     | 5       | U      |
| chloromethane            | 10 U                | 1,1,2-trichloroethane     | 5       | U      |
| dibromochloromethane     | 5 U                 | trichloroethene           | 5       | U      |
| 1,1-dichloroethane       | 5 U                 | trichlorofluoromethane    | - 5     | Ü      |
| 1,2-dichloroethane       | <b>5</b> , <b>U</b> | vinyl chloride            | 10      | U      |

Date of Analysis: 04/06/92

This method blank applies to the following sample: 5C-TF.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

lient Project ID: LE Carpenter

Job Number: ITEK 51037

# ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 5 Lab Sample ID: WB0406

Tentative Identification (1)

Concentration (2)

acetonitrile

54

Remarks:

(1) Identification is based on computer search of the NIST Library.

(2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITEK 51037

# WATER SURROGATE PERCENT RECOVERY SUMMARY

|                  |                       | VOLATILE          |                                     |
|------------------|-----------------------|-------------------|-------------------------------------|
| Client Sample ID | Toluene-D8 (88-110%)* | BFB<br>(86-115%)* | 1,2 Dichloroethane-D4<br>(76-114%)* |
| 1A-TF            | 105                   | 99                | 94                                  |
| 1B-TF            | 96                    | 95                | 94                                  |
| 1C-TF            | 103                   | 101               | 99                                  |
| 2A-TF            | 99                    | 97                | 96                                  |
| 2B-TF            | 95                    | 91                | 93                                  |
| 2C-TF            | 99                    | 96                | 95                                  |
| 3A-TF            | 97                    | 94                | 92                                  |
| 3B-TF            | 97                    | 92                | 89                                  |
| 3C-TF            | 104                   | 98                | 93                                  |
| 4A-TF            | 113] **               | 110               | 109                                 |
| 4A-TF DL         | 103                   | 102               | 81                                  |
| 4A-TF DL2        | 105                   | 110               | 105                                 |
| 4B-TF            | 103                   | 102               | 102                                 |
| 4B-TF DL         | 96                    | 96                | 91                                  |
| 4C-TF            | 101                   | 102               | 96                                  |
| 4C-TF DL         | 96                    | 99                | <b>98</b>                           |

<sup>\*</sup>Values in parenthesis represent QC limits.

DL - Dilution

<sup>\*\*</sup>Values are outside of QC limits.

Job Number: ITEK 51037

# WATER SURROGATE PERCENT RECOVERY SUMMARY

|                  | ****                     | VOLATILE          | •                                   |
|------------------|--------------------------|-------------------|-------------------------------------|
| Client Sample ID | Toluene-D8<br>(88-110%)* | BFB<br>(86-115%)* | 1,2 Dichloroethane-D4<br>(76-114%)* |
| 5A-TF            | 99                       | 102               | 94                                  |
| 5A-TF DL         | 89                       | 91                | 86                                  |
| 5B-TF            | 97                       | 101               | 93.                                 |
| 5B-TF DL         | 102                      | 102               | 95                                  |
| 5C-TF            | <b>95</b>                | 102               | 98                                  |
| Method Blank 1   | 99                       | 96                | 94                                  |
| Method Blank 2   | 108                      | 105               | 99                                  |
| Method Blank 3   | 100                      | 101               | 99                                  |
| Method Blank 4   | 94                       | 94                | 93                                  |
| Method Blank 5   | 99                       | 94                | 93                                  |

DL - Dilution

<sup>\*</sup>Values in parenthesis represent QC limits.

Job Number: ITEK 51037

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1A-TF Lab Sample ID: SS8682

| Compound                   | Concen | tration    | Compound                    | Concentr | ation |
|----------------------------|--------|------------|-----------------------------|----------|-------|
| acenaphthene               | 20     | U          | bis(2-chloroisopropyl)ether | 20       | U     |
| acenaphthylene             | 20     | U          | bis(2-ethylhexyl)phthalate  | 110      |       |
| anthracene                 | 20     | U          | 4-bromophenyl phenyl ether  | 20       | U     |
| benzidine                  | 100    | <b>u</b>   | 2-chloronaphthalene         | 20       | U     |
| benzo(a)anthracene         | 20     | Ü          | 4-chlorophenyl phenyl ether | 20       | U     |
| benzo(b)fluoranthene       | 20     | <b>U</b> , | chrysene                    | 20       | U     |
| benzo(k)fluoranthene       | 20     | U          | dibenz(a,h)anthracene       | 20       | U     |
| benzo(a)pyrene             | 20     | U          | di-n-butylphthalate         | 20       | U     |
| benzo(g,h,i)perylene       | 20     | U          | 1,2-dichlorobenzene         | 20       | U     |
| bütylbenzylphthalate       | 20     | U          | 1,3-dichlorobenzene         | 20       | U     |
| bis(2-chloroethoxy)methane | 20     | U          | 1,4-dichlorobenzene         | 20       | U     |
| bis(2-chloroethyl)ether    | 20     | U          |                             |          |       |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction:

04/03/92

Date of Analysis:

04/06/92

J - Indicates an estimated value less than the detection limit.

Job Number: ITEK 51037

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1A-TF Lab Sample ID: SS8682

| Compound                  | Concentration | Compound                   | Concentration |
|---------------------------|---------------|----------------------------|---------------|
| 3,3'dichlorobenzidine     | 40 U          | hexachloroethane           | 20 U          |
| diethylphthalate          | 20 U          | indeno(1,2,3-cd)pyrene     | 20 U          |
| dimethyl phthalate        | <b>20</b> U   | isophorone                 | <b>20</b> U   |
| 2,4-dinitrotoluene        | <b>20</b> U   | naphthalene                | <b>20</b> U   |
| 2,6-dinitrotoluene        | <b>20</b> U   | nitrobenzene               | 20 U          |
| di-n-octyl phthalate      | 20 Ü          | n-nitroso-di-n-propylamine | 20 U          |
| 1,2-diphenylhydrazine(1)  | 20 Ü          | n-nitrosodimethylamine     | 20 U          |
| fluoranthene              | 20 U          | n-nitrosodiphenylamine(2)  | <b>20</b> U   |
| fluorene                  | <b>20</b> U   | phenanthrene               | 20 U          |
| hexachlorobenzene         | <b>20</b> U   | pyrene                     | <b>20</b> U   |
| hexachlorobutadiene       | 20 U          | 1,2,4-trichlorobenzene     | <b>20</b> U   |
| hexachlorocyclopentadiene | <b>20</b> Ü   |                            |               |

- Compound was analyzed for but not detected. The number is the detection limit for the sample.

- Indicates an estimated value less than the detection limit.

(1) - Screened for as azobenzene

(2) - Detected as diphenylamine

Date of Extraction: Date of Analysis:

04/03/92 04/06/92

Client Project ID: LE Carpenter

Job Number: ITEK 51037

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in  $\mu g$ /liter (ppb)

Sample Matrix: Water

Client Sample ID: 1A-TF Lab Sample ID: SS8682

Tentative Identification (1)

Concentration (2)

NONE DETECTED

Remarks:

(1) Identification is based on computer search of the NIST Library.

(2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITEK 51037

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1B-TF Lab Sample ID: SS8683

| Compound                   | Concen | tration    | Compound                    | Concentr | ation |
|----------------------------|--------|------------|-----------------------------|----------|-------|
| acenaphthene               | 20     | U          | bis(2-chloroisopropyl)ether | 20       | Ü     |
| acenaphthylene             | 20     | U:         | bis(2-ethylhexyl)phthalate  | 200      |       |
| anthracene                 | 20     | U          | 4-bromophenyl phenyl ether  | 20       | U     |
| benzidine                  | 100    | Ŭ,         | 2-chloronaphthalene         | 20       | U     |
| benzo(a)anthracene         | 20     | U          | 4-chlorophenyl phenyl ether | 20       | U     |
| benzo(b)fluoranthene       | 20     | U          | chrysene                    | 20       | U     |
| benzo(k)fluoranthene       | 20     | U          | dibenz(a,h)anthracene       | 20       | U     |
| benzo(a)pyrene             | 20     | U          | di-n-butylphthalate         | 20       | U     |
| benzo(g,h,i)perylene       | 20     | . <b>U</b> | 1,2-dichlorobenzene         | 20       | U     |
| butylbenzylphthalate       | 20     | U          | 1,3-dichlorobenzene         | 20       | U     |
| bis(2-chloroethoxy)methane | 20     | U          | 1,4-dichlorobenzene         | 20       | Ų     |
| bis(2-chloroethyl)ether    | 20     | U          |                             |          |       |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction:

04/03/92

Date of Analysis:

04/06/92

J - Indicates an estimated value less than the detection limit.

Job Number: ITEK 51037

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1B-TF Lab Sample ID: SS8683

| Compound                  | Concentration | Compound                   | Concentration |
|---------------------------|---------------|----------------------------|---------------|
| 3,3'dichlorobenzidine     | 40 U          | hexachloroethane           | 20 U          |
| diethylphthalate          | <b>20</b> U   | indeno(1,2,3-cd)pyrene     | 20 U          |
| dimethyl phthalate        | 20 U          | isophorone                 | 20 U          |
| 2,4-dinitrotoluene        | 20 U          | naphthalene                | 20 U          |
| 2,6-dinitrotoluene        | <b>20</b> U   | nitrobenzene               | 20 U          |
| di-n-octyl phthalate      | <b>20</b> U   | n-nitroso-di-n-propylamine | 20 U          |
| 1,2-diphenylhydrazine(1)  | 20 U          | n-nitrosodimethylamine     | 20 U          |
| fluoranthene              | 20 U          | n-nitrosodiphenylamine(2)  | 20 U          |
| fluorene                  | 20 U          | phenanthrene               | 20 U          |
| hexachlorobenzene         | 20 U          | pyrene                     | 20 U          |
| hexachlorobutadiene       | 20 U          | 1,2,4-trichlorobenzene     | 20 U          |
| hexachlorocyclopentadiene | <b>20</b> U   |                            |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 04/03/92 Date of Analysis: 04/06/92

J - Indicates an estimated value less than the detection limit.

<sup>(1) -</sup> Screened for as azobenzene

<sup>(2) -</sup> Detected as diphenylamine

Client Project ID: LE Carpenter

Job Number: ITEK 51037

# ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1B-TF Lab Sample ID: SS8683

Tentative Identification (1)

NONE DETECTED

Concentration (2)

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITEK 51037

#### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1C-TF Lab Sample ID: SS8684

| Compound                   | Concentration | Compound                    | Concentration |  |
|----------------------------|---------------|-----------------------------|---------------|--|
| acenaphthene               | 20 U          | bis(2-chloroisopropyl)ether | 20 U          |  |
| acenaphthylene             | 20 U          | bis(2-ethylhexyl)phthalate  | 180           |  |
| anthracene                 | 20 U          | 4-bromophenyl phenyl ether  | 20 U          |  |
| benzidine                  | 100 U         | 2-chloronaphthalene         | .20 · U       |  |
| benzo(a)anthracene         | 20 U          | 4-chlorophenyl phenyl ether | 20 U          |  |
| benzo(b)fluoranthene       | 20 U          | chrysene                    | <b>20</b> U   |  |
| benzo(k)fluoranthene       | 20 U          | dibenz(a,h)anthracene       | 20 U          |  |
| benzo(a)pyrene             | 20 U          | di-n-butylphthalate         | <b>20</b> U   |  |
| benzo(g,h,i)perylene       | 20 U          | 1,2-dichlorobenzene         | ,20 U         |  |
| butylbenzylphthalate       | 20 U          | 1,3-dichlorobenzene         | 20 U          |  |
| bis(2-chloroethoxy)methane | 20 U          | 1,4-dichlorobenzene         | 20 U          |  |
| bis(2-chloroethyl)ether    | 20 U          |                             |               |  |

Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction:

04/03/92

Date of Analysis:

04/06/92

J - Indicates an estimated value less than the detection limit.

Job Number: ITEK 51037

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1C-TF Lab Sample ID: SS8684

| Compound                  | Concentration     | Compound                   | Concentration |
|---------------------------|-------------------|----------------------------|---------------|
| 3,3'dichlorobenzidine     | 40 U              | hexachloroethane           | 20 U          |
| diethylphthalate          | 20 U              | indeno(1,2,3-cd)pyrene     | 20 U          |
| dimethyl phthalate        | 20 U              | isophorone                 | 20 U          |
| 2,4-dinitrotoluene        | <b>20</b> U       | naphthalene                | 20 U          |
| 2,6-dinitrotoluene        | 20 U              | nitrobenzene               | <b>20</b> U   |
| di-n-octyl phthalate      | 20 U              | n-nitroso-di-n-propylamine | <b>20</b> U   |
| 1,2-diphenylhydrazine(1)  | 20 U              | n-nitrosodimethylamine     | 20 U          |
| fluoranthene              | <sup>"</sup> 20 U | n-nitrosodiphenylamine(2)  | 20 U          |
| fluorene                  | 20 U              | phenanthrene               | 20 U          |
| hexachlorobenzene         | 20 U              | pyrene                     | 20 U          |
| hexachlorobutadiene       | 20 U              | 1,2,4-trichlorobenzene     | 20 U          |
| hexachlorocyclopentadiene | 20 U              |                            |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

(1) - Screened for as azobenzene

(2) - Detected as diphenylamine

Date of Extraction:

04/03/92

Date of Analysis:

04/06/92

Client Project ID: LE Carpenter

Job Number: ITEK 51037

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1C-TF Lab Sample ID: SS8684

Tentative Identification (1)

Concentration (2)

NONE DETECTED

Remarks:

(1) Identification is based on computer search of the NIST Library.

(2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITEK 51037

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in  $\mu g$ /liter (ppb)

Sample Matrix: Water

Client Sample ID: 2A-TF Lab Sample ID: SS8685

| Compound                   | Concentration |     | Compound                    | Concentration |   |
|----------------------------|---------------|-----|-----------------------------|---------------|---|
| acenaphthene               | 20            | U , | bis(2-chloroisopropyl)ether | 20            | U |
| acenaphthylene             | 20            | U   | bis(2-ethylhexyl)phthalate  | 1,800         | D |
| anthracene                 | 20            | U   | 4-bromophenyl phenyl ether  | 20            | U |
| benzidine                  | 100           | U   | 2-chloronaphthalene         | 20            | U |
| benzo(a)anthracene         | 20            | U   | 4-chlorophenyl phenyl ether | 20            | U |
| benzo(b)fluoranthene       | 20            | U   | chrysene                    | 20            | Ü |
| benzo(k)fluoranthene       | 20            | U · | dibenz(a,h)anthracene       | 20            | U |
| benzo(a)pyrene             | 20            | U.  | di-n-butylphthalate         | 20            | U |
| benzo(g,h,i)perylene       | 20            | U   | 1,2-dichlorobenzene         | 20            | U |
| butylbenzylphthalate       | 6             | J   | 1,3-dichlorobenzene         | 20            | U |
| bis(2-chloroethoxy)methane | 20            | U   | 1,4-dichlorobenzene         | 20            | U |
| bis(2-chloroethyl)ether    | 20            | U   |                             |               |   |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 04/03/92 Date of Analysis: 04/06/92

J - Indicates an estimated value less than the detection limit.

D - Compound analyzed at a secondary dilution factor.

Job Number: ITEK 51037

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 2A-TF Lab Sample ID: SS8685

| Compound                  | Concentration | Compound                   | Concentration |  |  |
|---------------------------|---------------|----------------------------|---------------|--|--|
| 3,3'dichlorobenzidine     | 40 U          | hexachloroethane           | 20 Ü          |  |  |
| diethylphthalate          | 20 U          | indeno(1,2,3-cd)pyrene     | 20 · U        |  |  |
| dimethyl phthalate        | <b>20</b> U   | isophorone                 | 20 U          |  |  |
| 2,4-dinitrotoluene        | <b>20</b> U   | naphthalene                | 20 U          |  |  |
| 2,6-dinitrotoluene        | 20 U          | nitrobenzene               | 20 U          |  |  |
| di-n-octyl phthalate      | <b>20</b> U   | n-nitroso-di-n-propylamine | 20 U          |  |  |
| 1,2-diphenylhydrazine(1)  | <b>20</b> U   | n-nitrosodimethylamine     | 20 Ü          |  |  |
| fluoranthene              | <b>20</b> U   | n-nitrosodiphenylamine(2)  | 20 U          |  |  |
| fluorene                  | 20 U          | phenanthrene               | 20 U          |  |  |
| hexachlorobenzene         | <b>20 U</b>   | pyrene                     | 20 U          |  |  |
| hexachlorobutadiene       | 20 U          | 1,2,4-trichlorobenzene     | 20 U          |  |  |
| hexachlorocyclopentadiene | <b>20</b> U   |                            | •             |  |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Indicates an estimated value less than the detection limit.

(1) - Screened for as azobenzene

(2) - Detected as diphenylamine

Date of Extraction:

04/03/92

Date of Analysis:

04/06/92

Client Project ID: LE Carpenter

Job Number: ITEK 51037

# ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 2A-TF Lab Sample ID: SS8685

| Tentative Identification (1)            | Concentration | 1 (2) |
|---|---------------|-------|
| 2-butanone, 4-(dimethylamino)-3-methyl- | 23            | J     |
| unknown                                 | 35            | J ·   |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- J Indicates an estimated value less than the detection limit.

Job Number: ITEK 51037

#### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 2B-TF Lab Sample ID: SS8686

| Compound                   | Concentration |     | Compound                    | Concentration |   |
|----------------------------|---------------|-----|-----------------------------|---------------|---|
| acenaphthene               | 20            | U   | bis(2-chloroisopropyl)ether | 20            | U |
| acenaphthylene             | 20            | U   | bis(2-ethylhexyl)phthalate  | 1,000         | D |
| anthracene                 | 20            | U   | 4-bromophenyl phenyl ether  | 20            | U |
| benzidine                  | 100           | Ü   | 2-chloronaphthalene         | 20            | U |
| benzo(a)anthracene         | 20            | U   | 4-chlorophenyl phenyl ether | 20            | U |
| benzo(b)fluoranthene       | 20            | U   | chrysene                    | 20            | U |
| benzo(k)fluoranthene       | 20            | U   | dibenz(a,h)anthracene       | 20            | U |
| benzo(a)pyrene             | 20            | U - | di-n-butylphthalate         | 20            | U |
| benzo(g,h,i)perylene       | 20            | U - | 1,2-dichlorobenzene         | 20            | U |
| butylbenzylphthalate       | 20            | U   | 1,3-dichlorobenzene         | 20            | U |
| bis(2-chloroethoxy)methane | 20            | U   | 1,4-dichlorobenzene         | 20            | U |
| bis(2-chloroethyl)ether    | 20            | U   |                             |               |   |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction:

04/03/92

Date of Analysis:

04/07/92

J - Indicates an estimated value less than the detection limit.

D - Compound analyzed at a secondary dilution factor.

Job Number: ITEK 51037

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 2B-TF Lab Sample ID: SS8686

| Compound                  | Concentration | Compound                   | Concentration |  |
|---------------------------|---------------|----------------------------|---------------|--|
| 3,3'dichlorobenzidine     | 40 U          | hexachloroethane           | 20 Ų          |  |
| diethylphthalate          | 20 U          | indeno(1,2,3-cd)pyrene     | 20 U          |  |
| dimethyl phthalate        | 20 U          | isophorone                 | 20 U          |  |
| 2,4-dinitrotoluene        | 20 U          | naphthalene                | 20 U          |  |
| 2,6-dinitrotoluene        | 20 Ü          | nitrobenzene               | 20 U          |  |
| di-n-octyl phthalate      | 20 U          | n-nitroso-di-n-propylamine | 20 U          |  |
| 1,2-diphenylhydrazine(1)  | 20 U          | n-nitrosodimethylamine     | 20 U          |  |
| fluoranthene              | 20 U          | n-nitrosodiphenylamine(2)  | 20 U          |  |
| fluorene                  | 20 U          | phenanthrene               | 20 U          |  |
| hexachlorobenzene         | <b>20</b> U   | pyrene                     | 20 U          |  |
| hexachlorobutadiene       | 20 Ü          | 1,2,4-trichlorobenzene     | <b>20</b> U   |  |
| hexachlorocyclopentadiene | 20 U          |                            |               |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

(1) - Screened for as azobenzene

(2) - Detected as diphenylamine

Date of Extraction:

04/03/92

Date of Analysis:

04/07/92

Client Project ID: LE Carpenter

Job Number: ITEK 51037

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in  $\mu g/\text{liter}$  (ppb)

Sample Matrix: Water

Client Sample ID: 2B-TF Lab Sample ID: SS8686

| Tentative Identification (1)            | Concentration | <u>(2)</u> |
|---|---------------|------------|
| 2-butanone, 4-(dimethylamino)-3-methyl- | 18            | J          |
| unknown                                 | 25            | J          |
| unknown                                 | 24            | J          |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- J Indicates an estimated value less than the detection limit.

Job Number: ITEK 51037

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 2C-TF Lab Sample ID: SS8687

| Compound                   | Concentration |                   | Compound                    | Concentration |   |
|----------------------------|---------------|-------------------|-----------------------------|---------------|---|
| acenaphthene               | 20            | <b>U</b>          | bis(2-chloroisopropyl)ether | 20            | U |
| acenaphthylene             | 20            | U                 | bis(2-ethylhexyl)phthalate  | 1,900         | D |
| anthracene                 | 20            | U -               | 4-bromophenyl phenyl ether  | 20            | Ü |
| benzidine                  | 100           | U ,               | 2-chloronaphthalene         | 20            | U |
| benzo(a)anthracene         | 20            | U                 | 4-chlorophenyl phenyl ether | 20            | U |
| benzo(b)fluoranthene       | 20            | U                 | chrysene                    | 20            | U |
| benzo(k)fluoranthene       | 20            | U                 | dibenz(a,h)anthracene       | 20            | U |
| benzo(a)pyrene             | 20            | U                 | di-n-butylphthalate         | 20            | Ų |
| benzo(g,h,i)perylene       | 20            | $\mathbf{U}^{-1}$ | 1,2-dichlorobenzene         | 20            | U |
| butylbenzylphthalate       | 4             | J                 | 1,3-dichlorobenzene         | 20            | U |
| bis(2-chloroethoxy)methane | 20            | U                 | 1,4-dichlorobenzene         | 20            | U |
| bis(2-chloroethyl)ether    | 20            | U                 |                             |               |   |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 04/03/92 Date of Analysis: 04/07/92

J - Indicates an estimated value less than the detection limit.

D - Compound analyzed at a secondary dilution factor.

Job Number: ITEK 51037

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 2C-TF Lab Sample ID: SS8687

| Compound                  | Concentration |           | Compound                   | Concentration |   |  |
|---------------------------|---------------|-----------|----------------------------|---------------|---|--|
| 3,3'dichlorobenzidine     | 40            | U         | hexachloroethane           | 20            | Ų |  |
| diethylphthalate          | 20            | U         | indeno(1,2,3-cd)pyrene     | 20            | U |  |
| dimethyl phthalate        | 20            | U         | isophorone                 | 20            | U |  |
| 2,4-dinitrotoluene        | 20            | <b>U</b>  | naphthalene                | 20            | U |  |
| 2,6-dinitrotoluene        | 20            | U         | nitrobenzene               | 20            | U |  |
| di-n-octyl phthalate      | 20            | U         | n-nitroso-di-n-propylamine | 20            | U |  |
| 1,2-diphenylhydrazine(1)  | 20            | U.        | n-nitrosodimethylamine     | 20            | U |  |
| fluoranthene              | 20            | U         | n-nitrosodiphenylamine(2)  | 20            | U |  |
| fluorene                  | 20            | U         | phenanthrene               | 20            | U |  |
| hexachlorobenzene         | 20            | <b>"U</b> | pyrene                     | 20            | U |  |
| hexachlorobutadiene       | 20            | U         | 1,2,4-trichlorobenzene     | 20            | U |  |
| hexachlorocyclopentadiene | 20            | U         |                            |               | • |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Indicates an estimated value less than the detection limit.

(1) - Screened for as azobenzene

(2) - Detected as diphenylamine

Date of Extraction: 0
Date of Analysis: 0

04/03/92

04/07/92

Client Project ID: LE Carpenter

Job Number: ITEK 51037

# ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 2C-TF Lab Sample ID: SS8687

| Tentative Identification (1)            | Concentration (2) |  |
|---|-------------------|--|
| 2-butanone, 4-(dimethylamino)-3-methyl- | 26 J              |  |
| unknown                                 | 40 J              |  |
| unknown                                 | 18 J              |  |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- J Indicates an estimated value less than the detection limit.

Job Number: ITEK 51037

### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 3A-TF Lab Sample ID: SS8688

| Compound                   | Concentration |              | Compound                    | Concentr | <b>Concentration</b> |  |
|----------------------------|---------------|--------------|-----------------------------|----------|----------------------|--|
| acenaphthene               | 20            | U            | bis(2-chloroisopropyl)ether | 20       | U                    |  |
| acenaphthylene             | 20            | U            | bis(2-ethylhexyl)phthalate  | 150      |                      |  |
| anthracene                 | 20            | U            | 4-bromophenyl phenyl ether  | 20       | U                    |  |
| benzidine                  | 100           | U            | 2-chloronaphthalene         | 20       | U                    |  |
| benzo(a)anthracene         | 20            | <b>U</b>     | 4-chlorophenyl phenyl ether | 20       | U                    |  |
| benzo(b)fluoranthene       | 20            | U            | chrysene                    | 20       | Ü                    |  |
| benzo(k)fluoranthene       | 20            | U            | dibenz(a,h)anthracene       | 20       | U                    |  |
| benzo(a)pyrene             | 20            | <b>U</b> .,, | di-n-butylphthalate         | 20       | U                    |  |
| benzo(g,h,i)perylene       | 20            | U            | 1,2-dichlorobenzene         | 20       | U                    |  |
| butylbenzylphthalate       | 20            | U            | 1,3-dichlorobenzene         | 20       | Ü                    |  |
| bis(2-chloroethoxy)methane | 20            | U            | 1,4-dichlorobenzene         | 20       | Ü                    |  |
| bis(2-chloroethyl)ether    | 20            | U            |                             |          |                      |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction:

04/03/92

Date of Analysis:

04/07/92

J - Indicates an estimated value less than the detection limit.

Job Number: ITEK 51037

## BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 3A-TF Lab Sample ID: SS8688

| Compound                  | Concentra | ation | Compound                   | Concentra | ation |
|---------------------------|-----------|-------|----------------------------|-----------|-------|
| 3,3'dichlorobenzidine     | 40        | U     | hexachloroethane           | 20        | Ų     |
| diethylphthalate          | 20        | Ü     | indeno(1,2,3-cd)pyrene     | 20        | U     |
| dimethyl phthalate        | 20        | Ü     | isophorone                 | 20        | Ų     |
| 2,4-dinitrotoluene        | 20        | U     | naphthalene                | 20        | U     |
| 2,6-dinitrotoluene        | 20        | U     | nitrobenzene               | 20        | U     |
| di-n-octyl phthalate      | 20        | Ŭ     | n-nitroso-di-n-propylamine | 20        | U     |
| 1,2-diphenylhydrazine(1)  | 20        | Ü     | n-nitrosodimethylamine     | 20        | U     |
| fluoranthene              | 20        | U     | n-nitrosodiphenylamine(2)  | 20        | Ŭ     |
| fluorene                  | 20        | U     | phenanthrene               | 20        | U     |
| hexachlorobenzene         | 20        | U     | pyrene                     | 20        | U     |
| hexachlorobutadiene       | 20        | U     | 1,2,4-trichlorobenzene     | 20        | U     |
| hexachlorocyclopentadiene | 20        | U     |                            |           |       |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

(1) - Screened for as azobenzene

(2) - Detected as diphenylamine

Date of Extraction:
Date of Analysis:

04/03/92

04/07/92

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

Client Project ID: LE Carpenter

Job Number: ITEK 51037

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 3A-TF Lab Sample ID: SS8688

| Tentative Identification (1)     |       | Concentration | (2) |
|----------------------------------|-------|---------------|-----|
| 2-pentanone, 4-hydroxy-4-methyl- |       | 17            | JAB |
| unknown                          | 1 - 1 | 21            | J   |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- A Suspected aldol product.
- B Analyte was found in the blank as well as the sample.
- J Indicates an estimated value less than the detection limit.

Job Number: ITEK 51037

#### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in  $\mu g$ /liter (ppb)

Sample Matrix: Water

Client Sample ID: 3B-TF Lab Sample ID: SS8689

| Compound                   | Concen | tration | Compound                    | Concentr | ation |
|----------------------------|--------|---------|-----------------------------|----------|-------|
| acenaphthene               | 20     | Ų       | bis(2-chloroisopropyl)ether | 20       | U     |
| acenaphthylene             | 20     | U       | bis(2-ethylhexyl)phthalate  | 86       |       |
| anthracene                 | 20     | U .     | 4-bromophenyl phenyl ether  | 20       | U     |
| benzidine                  | 100    | U       | 2-chloronaphthalene         | 20       | U     |
| benzo(a)anthracene         | 20     | Ü       | 4-chlorophenyl phenyl ether | 20       | U     |
| benzo(b)fluoranthene       | 20     | U       | chrysene                    | 20       | U     |
| benzo(k)fluoranthene       | 20     | U       | dibenz(a,h)anthracene       | 20       | Ų     |
| benzo(a)pyrene             | 20     | U       | di-n-butylphthalate         | 20       | U     |
| benzo(g,h,i)perylene       | 20     | U .     | 1,2-dichlorobenzene         | 20       | U     |
| butylbenzylphthalate       | 20     | U       | 1,3-dichlorobenzene         | .20      | U     |
| bis(2-chloroethoxy)methane | 20     | U ·     | 1,4-dichlorobenzene         | .20      | U     |
| bis(2-chloroethyl)ether    | 20     | U       |                             |          |       |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction:

04/03/92

Date of Analysis:

04/18/92

J - Indicates an estimated value less than the detection limit.

Job Number: ITEK 51037

### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in  $\mu g$ /liter (ppb)

Sample Matrix: Water

Client Sample ID: 3B-TF Lab Sample ID: SS8689

| Compound                  | <b>Concentration</b> | Compound                   | <b>Concentration</b> |
|---------------------------|----------------------|----------------------------|----------------------|
| 3,3'dichlorobenzidine     | 40 U                 | hexachloroethane           | 20 U                 |
| diethylphthalate          | 20 U                 | indeno(1,2,3-cd)pyrene     | 20 U                 |
| dimethyl phthalate        | 20 U                 | isophorone                 | . 20 U               |
| 2,4-dinitrotoluene        | 20 U                 | naphthalene                | 20 U                 |
| 2,6-dinitrotoluene        | 20 Ü                 | nitrobenzene               | .20 U                |
| di-n-octyl phthalate      | <b>20</b> U          | n-nitroso-di-n-propylamine | 20 U                 |
| 1,2-diphenylhydrazine(1)  | 20 U                 | n-nitrosodimethylamine     | <b>20</b> U          |
| fluoranthene              | 20 U                 | n-nitrosodiphenylamine(2)  | 20 U                 |
| fluorene                  | 20 U                 | phenanthrene               | 20 U                 |
| hexachlorobenzene         | 20 U                 | pyrene                     | 20 U                 |
| hexachlorobutadiene       | 20 U                 | 1,2,4-trichlorobenzene     | 20 U                 |
| hexachlorocyclopentadiene | 20 U                 |                            |                      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

- Indicates an estimated value less than the detection limit.

(1) - Screened for as azobenzene

(2) - Detected as diphenylamine

Date of Extraction: 04/03/92 Date of Analysis: 04/18/92

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

Client Project ID: LE Carpenter

Job Number: ITEK 51037

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 3B-TF Lab Sample ID: SS8689

Tentative Identification (1)

Concentration (2)

NONE DETECTED

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITEK 51037

#### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in  $\mu g/liter$  (ppb)

Sample Matrix: Water

Client Sample ID: 3C-TF Lab Sample ID: SS8690

| Compound                   | Concen | tration    | Compound                    | Concentr | ration |
|----------------------------|--------|------------|-----------------------------|----------|--------|
| acenaphthene               | 20     | U          | bis(2-chloroisopropyl)ether | 20       | U      |
| acenaphthylene             | 20     | U          | bis(2-ethylhexyl)phthalate  | . 27     |        |
| anthracene                 | 20     | U,         | 4-bromophenyl phenyl ether  | 20       | U      |
| benzidine                  | 100    | U          | 2-chloronaphthalene         | 20       | U      |
| benzo(a)anthracene         | 20     | U          | 4-chlorophenyl phenyl ether | 20       | U      |
| benzo(b)fluoranthene       | 20     | U ,        | chrysene                    | 20       | U      |
| benzo(k)fluoranthene       | 20     | , <b>U</b> | dibenz(a,h)anthracene       | 20       | U      |
| benzo(a)pyrene             | 20     | U ·        | di-n-butylphthalate         | 20       | Ù      |
| benzo(g,h,i)perylene       | 20     | U          | 1,2-dichlorobenzene         | 20       | U      |
| butylbenzylphthalate       | 20     | U          | 1,3-dichlorobenzene         | 20       | U      |
| bis(2-chloroethoxy)methane | 20     | U          | 1,4-dichlorobenzene         | 20       | U      |
| bis(2-chloroethyl)ether    | 20     | U          |                             | •        |        |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction:

04/03/92

Date of Analysis:

04/18/92

J - Indicates an estimated value less than the detection limit.

Job Number: ITEK 51037

### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 3C-TF Lab Sample ID: SS8690

| Compound                  | Concentration | Compound                   | Concentration |
|---------------------------|---------------|----------------------------|---------------|
| 3,3'dichlorobenzidine     | 40 U          | hexachloroethane           | <b>20</b> Ų   |
| diethylphthalate          | 20 Ú          | indeno(1,2,3-cd)pyrene     | <b>20</b> U   |
| dimethyl phthalate        | 20 U          | isophorone                 | 20 U          |
| 2,4-dinitrotoluene        | 20 U          | naphthalene                | 20 U          |
| 2,6-dinitrotoluene        | 20 U          | nitrobenzene               | 20 U          |
| di-n-octyl phthalate      | 20 U          | n-nitroso-di-n-propylamine | 20 U          |
| 1,2-diphenylhydrazine(1)  | 20 U          | n-nitrosodimethylamine     | 20 U          |
| fluoranthene              | 20 U          | n-nitrosodiphenylamine(2)  | 20 U          |
| fluorene                  | 20 U          | phenanthrene               | 20 U          |
| hexachlorobenzene         | 20 U          | pyrene                     | <b>20</b> U   |
| hexachlorobutadiene       | 20 U          | 1,2,4-trichlorobenzene     | 20 U          |
| hexachlorocyclopentadiene | 20 U          |                            |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

(1) - Screened for as azobenzene

(2) - Detected as diphenylamine

Date of Extraction:

04/03/92

Date of Analysis:

04/18/92

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

'lient Project ID: LE Carpenter

Job Number: ITEK 51037

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 3C-TF Lab Sample ID: SS8690

Tentative Identification (1)

Concentration (2)

NONE DETECTED

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITEK 51037

#### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in  $\mu g/liter$  (ppb)

Sample Matrix: Water

Client Sample ID: 4A-TF Lab Sample ID: SS8691

| Compound                   | Concent | ration       | Compound                    | Concenti | ration |
|----------------------------|---------|--------------|-----------------------------|----------|--------|
| acenaphthene               | 20      | U            | bis(2-chloroisopropyl)ether | 20       | U      |
| acenaphthylene             | 20      | U            | bis(2-ethylhexyl)phthalate  | 11,000   | D      |
| anthracene                 | 20      | U            | 4-bromophenyl phenyl ether  | 20       | U      |
| benzidine                  | 100     | U            | 2-chloronaphthalene         | 20       | U      |
| benzo(a)anthracene         | 20      | U            | 4-chlorophenyl phenyl ether | 20       | U      |
| benzo(b)fluoranthene       | 20      | U            | chrysene                    | 20       | Ü      |
| benzo(k)fluoranthene       | 20      | U            | dibenz(a,h)anthracene       | 20       | Ü      |
| benzo(a)pyrene             | 20      | U'.          | di-n-butylphthalate         | 20       | U      |
| benzo(g,h,i)perylene       | 20      | <b>U</b> , i | 1,2-dichlorobenzene         | 20       | U      |
| butylbenzylphthalate       | 20      | U            | 1,3-dichlorobenzene         | 20       | U      |
| bis(2-chloroethoxy)methane | 20      | U            | 1,4-dichlorobenzene         | 20       | U      |
| bis(2-chloroethyl)ether    | 20      | Ü,           |                             |          |        |

Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction:

04/03/92

Date of Analysis:

04/18/92

Indicates an estimated value less than the detection limit.

Compound analyzed at a secondary dilution factor.

KNOXVILLE, TN

Job Number: ITEK 51037

#### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 4A-TF Lab Sample ID: SS8691

| Compound                  | Concentra | ation      | Compound                   | Concentra | ation |
|---------------------------|-----------|------------|----------------------------|-----------|-------|
| 3,3'dichlorobenzidine     | 40        | U          | hexachloroethane           | 20        | Ų     |
| diethylphthalate          | 20        | Ü          | indeno(1,2,3-cd)pyrene     | 20        | U     |
| dimethyl phthalate        | 20        | Ü          | isophorone                 | .20       | U     |
| 2,4-dinitrotoluene        | 20        | Ü          | naphthalene                | 20        | U     |
| 2,6-dinitrotoluene        | 20        | U          | nitrobenzene               | 20        | U     |
| di-n-octyl phthalate      | 4         | J          | n-nitroso-di-n-propylamine | 20        | U     |
| 1,2-diphenylhydrazine(1)  | 20        | U          | n-nitrosodimethylamine     | .20       | U     |
| fluoranthene              | 20        | U          | n-nitrosodiphenylamine(2)  | .20       | U     |
| fluorene                  | 20        | Ü          | phenanthrene               | 20        | U     |
| hexachlorobenzene         | 20        | <b>U</b> : | pyrene                     | 20        | U     |
| hexachlorobutadiene       | 20        | U          | 1,2,4-trichlorobenzene     | 20        | U.    |
| hexachlorocyclopentadiene | 20        | U          |                            |           |       |

- Compound was analyzed for but not detected. The number is the detection limit for the sample.

- Indicates an estimated value less than the detection limit.

Screened for as azobenzene

(2) - Detected as diphenylamine

Date of Extraction: Date of Analysis:

04/03/92 04/18/92

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

Client Project ID: LE Carpenter

Job Number: ITEK 51037

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 4A-TF Lab Sample ID: SS8691

| Tentative Identification (1)   | Concentration ( | <u>2)</u> |
|--------------------------------|-----------------|-----------|
| benzene, (1-methylethyl)-      | 20 J            |           |
| benzaldehyde                   | 32 J            |           |
| benzenemethanol, .alphamethyl- | 24 J            |           |
| 2-propenal, 3-phenyl-          | 24 J            |           |
| unknown (phthalate)            | 48 J            |           |

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

J - Indicates an estimated value less than the detection limit.

Job Number: ITEK 51037

#### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 4B-TF Lab Sample ID: SS8692

| Compound                   | Concen | tration | Compound                    | Concentr | ation    |
|----------------------------|--------|---------|-----------------------------|----------|----------|
| acenaphthene               | 20     | U       | bis(2-chloroisopropyl)ether | 20       | U        |
| acenaphthylene             | 20     | U       | bis(2-ethylhexyl)phthalate  | 3,400    | D        |
| anthracene                 | 20     | U       | 4-bromophenyl phenyl ether  | 20       | U        |
| benzidine                  | 100    | U .     | 2-chloronaphthalene         | 20       | U        |
| benzo(a)anthracene         | 20     | U       | 4-chlorophenyl phenyl ether | 20       | U        |
| benzo(b)fluoranthene       | 20     | U ,     | chrysene                    | 20       | U        |
| benzo(k)fluoranthene       | 20     | U       | dibenz(a,h)anthracene       | 20       | U        |
| benzo(a)pyrene             | 20     | Ŭ,      | di-n-butylphthalate         | 20       | U        |
| benzo(g,h,i)perylene       | 20     | U       | 1,2-dichlorobenzene         | 20       | <b>U</b> |
| butylbenzylphthalate       | 20     | U       | 1,3-dichlorobenzene         | 20       | U        |
| bis(2-chloroethoxy)methane | 20     | U       | 1,4-dichlorobenzene         | 20       | U        |
| bis(2-chloroethyl)ether    | 20     | U       |                             |          |          |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 04/03/92 Date of Analysis: 04/18/92

J - Indicates an estimated value less than the detection limit.

D - Compound analyzed at a secondary dilution factor.

Job Number: ITEK 51037

### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 4B-TF Lab Sample ID: SS8692

| Compound                  | Concentra | tion | Compound                   | Concentra | tion |
|---------------------------|-----------|------|----------------------------|-----------|------|
| 3,3'dichlorobenzidine     | 40        | U    | hexachloroethane           | 20        | Ų    |
| diethylphthalate          | 20        | U-   | indeno(1,2,3-cd)pyrene     | 20        | U    |
| dimethyl phthalate        | 20        | U    | isophorone                 | 20        | U    |
| 2,4-dinitrotoluene        | 20        | U    | naphthalene                | 20        | U    |
| 2,6-dinitrotoluene        | 20        | Ŭ    | nitrobenzene               | 20        | U    |
| di-n-octyl phthalate      | 20        | Ū    | n-nitroso-di-n-propylamine | 20        | Ū    |
| 1,2-diphenylhydrazine(1)  | 20        | U    | n-nitrosodimethylamine     | 20        | Ü    |
| fluoranthene              | 20        | U    | n-nitrosodiphenylamine(2)  | 20        | U    |
| fluorene                  | 20        | U    | phenanthrene               | 20        | U    |
| hexachlorobenzene         | 20        | Ų    | pyrene                     | 20        | U    |
| hexachlorobutadiene       | 20        | U    | 1,2,4-trichlorobenzene     | 20        | U    |
| hexachlorocyclopentadiene | 20        | Ù    |                            |           |      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

(1) - Screened for as azobenzene

(2) - Detected as diphenylamine

Date of Extraction: 04/03/92 Date of Analysis: 04/18/92

Job Number: ITEK 51037

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 4B-TF Lab Sample ID: SS8692

| Tentative Identification (1)   | Concentration (2) |    |  |
|--------------------------------|-------------------|----|--|
| benzene, (1-methylethyl)-      | 21                | J  |  |
| benzaldehyde                   | 49                | J  |  |
| benzenemethanol, .alphamethyl- | 36                | J  |  |
| methyl benzaldehyde            | 22                | JY |  |
| 2-propenal, 3-phenyl-          | 31                | J  |  |
| unknown (phthalate)            | 21                | J  |  |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- Indicates an estimated value less than the detection limit.
- Y Indistinguishable isomer in tentatively identified compounds.

Job Number: ITEK 51037

#### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 4C-TF Lab Sample ID: SS8693

| Compound                   | Concentration | Compound                    | Concentration |
|----------------------------|---------------|-----------------------------|---------------|
| acenaphthene               | <b>20</b> U   | bis(2-chloroisopropyl)ether | <b>20</b> U   |
| acenaphthylene             | 20 U          | bis(2-ethylhexyl)phthalate  | 3,400 D       |
| anthracene                 | 20 U          | 4-bromophenyl phenyl ether  | 20 U          |
| benzidine                  | 100 U         | 2-chloronaphthalene         | 20 U          |
| benzo(a)anthracene         | 20 Ŭ          | 4-chlorophenyl phenyl ether | 20 U          |
| benzo(b)fluoranthene       | 20 U          | chrysene                    | <b>20</b> U   |
| benzo(k)fluoranthene       | 20 U          | dibenz(a,h)anthracene       | <b>20</b> U   |
| benzo(a)pyrene             | 20 U          | di-n-butylphthalate         | <b>20</b> U   |
| benzo(g,h,i)perylene       | 20 U          | 1,2-dichlorobenzene         | 20 U          |
| butylbenzylphthalate       | 20 U          | 1,3-dichlorobenzene         | 20 U          |
| bis(2-chloroethoxy)methane | 20 U          | 1,4-dichlorobenzene         | 20 U          |
| bis(2-chloroethyl)ether    | <b>20</b> U   |                             |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction:

04/03/92

Date of Analysis:

04/18/92

J - Indicates an estimated value less than the detection limit.

D - Compound analyzed at a secondary dilution factor.

Job Number: ITEK 51037

### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 4C-TF Lab Sample ID: SS8693

| Compound                  | Concentration | Compound                   | Concentration |
|---------------------------|---------------|----------------------------|---------------|
| 3,3'dichlorobenzidine     | <b>40</b> U   | hexachloroethane           | 20 U          |
| diethylphthalate          | 20 U          | indeno(1,2,3-cd)pyrene     | 20 U          |
| dimethyl phthalate        | 20 U          | isophorone                 | <b>20</b> U   |
| 2,4-dinitrotoluene        | <b>20</b> U   | naphthalene                | 20 U          |
| 2,6-dinitrotoluene        | 20 Ü          | nitrobenzene               | 20 U          |
| di-n-octyl phthalate      | 20 U          | n-nitroso-di-n-propylamine | 20 U          |
| 1,2-diphenylhydrazine(1)  | 20 U          | n-nitrosodimethylamine     | <b>20</b> U   |
| fluoranthene              | 20 U          | n-nitrosodiphenylamine(2)  | <b>20</b> U   |
| fluorene                  | 20 U          | phenanthrene               | 20 U          |
| hexachlorobenzene         | 20 U          | pyrene                     | 20 U          |
| hexachlorobutadiene       | 20 U          | 1,2,4-trichlorobenzene     | 20 U          |
| hexachlorocyclopentadiene | 20 U          | •                          |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 04/03/92 Date of Analysis: 04/18/92

J - Indicates an estimated value less than the detection limit.

<sup>(1) -</sup> Screened for as azobenzene

<sup>(2) -</sup> Detected as diphenylamine

Job Number: ITEK 51037

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 4C-TF Lab Sample ID: SS8693

| Tentative Identification (1)     | Concentration (2) |    |  |
|----------------------------------|-------------------|----|--|
| benzene, (1-methylethyl)-        | 21                | J  |  |
| benzaldehyde                     | 60                | J  |  |
| benzenemethanol, .alpha,-methyl- | 28                | J  |  |
| methyl benzaldehyde              | 41                | JY |  |
| 2-propenal, 3-phenyl-            | 65                | J  |  |
| unknown (phthalate)              | 21                | J  |  |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- J Indicates an estimated value less than the detection limit.
- Y Indistinguishable isomer in tentatively identified compounds.

Job Number: ITEK 51037

#### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 5A-TF Lab Sample ID: SS8694

| Compound                   | Concentration | Compound                    | Concentration |  |  |
|----------------------------|---------------|-----------------------------|---------------|--|--|
| acenaphthene               | <b>2</b> 0 U  | bis(2-chloroisopropyl)ether | 20 U          |  |  |
| acenaphthylene             | <b>20</b> U   | bis(2-ethylhexyl)phthalate  | 4,200 D       |  |  |
| anthracene                 | 20 U          | 4-bromophenyl phenyl ether  | 20 U          |  |  |
| benzidine                  | 1 <b>00</b> U | 2-chloronaphthalene         | 20 U          |  |  |
| benzo(a)anthracene         | <b>20</b> U   | 4-chlorophenyl phenyl ether | 20 U          |  |  |
| benzo(b)fluoranthene       | <b>20</b> U   | chrysene                    | 20 U          |  |  |
| benzo(k)fluoranthene       | <b>20</b> U   | dibenz(a,h)anthracene       | 20 U          |  |  |
| benzo(a)pyrene             | 20 U          | di-n-butylphthalate         | 20 U          |  |  |
| benzo(g,h,i)perylene       | 20 U          | 1,2-dichlorobenzene         | <b>20</b> U   |  |  |
| butylbenzylphthalate       | <b>2</b> 0 U  | 1,3-dichlorobenzene         | 20 U          |  |  |
| bis(2-chloroethoxy)methane | 20 U          | 1,4-dichlorobenzene         | 20 U          |  |  |
| bis(2-chloroethyl)ether    | 20 U          |                             |               |  |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 04/03/92 Date of Analysis: 04/18/92

J - Indicates an estimated value less than the detection limit.

D - Compound analyzed at a secondary dilution factor.

Job Number: ITEK 51037

### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 5A-TF Lab Sample ID: SS8694

| Compound                  | Concentration | Compound                   | Concentration |
|---------------------------|---------------|----------------------------|---------------|
| 3,3'dichlorobenzidine     | 40 U          | hexachloroethane           | 20 Ų          |
| diethylphthalate          | 20 U          | indeno(1,2,3-cd)pyrene     | 20 U          |
| dimethyl phthalate        | 20 U          | isophorone                 | 20 U          |
| 2,4-dinitrotoluene        | 20 U          | naphthalene                | 20 U          |
| 2,6-dinitrotoluene        | 20 U          | nitrobenzene               | 20 U          |
| di-n-octyl phthalate      | 20 U          | n-nitroso-di-n-propylamine | 20 U          |
| 1,2-diphenylhydrazine(1)  | 20 U          | n-nitrosodimethylamine     | <b>20</b> U   |
| fluoranthene              | 20 U          | n-nitrosodiphenylamine(2)  | <b>20</b> U   |
| fluorene                  | 20 U          | phenanthrene               | 20 U          |
| hexachlorobenzene         | 20 U          | pyrene                     | <b>20</b> U   |
| hexachlorobutadiene       | 20 U          | 1,2,4-trichlorobenzene     | 20 U          |
| hexachlorocyclopentadiene | 20 U          |                            |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 04/03/92 Date of Analysis: 04/18/92

J - Indicates an estimated value less than the detection limit.

<sup>(1) -</sup> Screened for as azobenzene

<sup>(2) -</sup> Detected as diphenylamine

Job Number: ITEK 51037

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in  $\mu g/\text{liter (ppb)}$ 

Sample Matrix: Water

Client Sample ID: 5A-TF Lab Sample ID: SS8694

| Tentative Identification (1)   | Concentration (2) |  |  |  |
|--------------------------------|-------------------|--|--|--|
| benzene, (1-methylethyl)-      | 46 J              |  |  |  |
| benzenemethanol, .alphamethyl- | 33 J              |  |  |  |
| ethanone, 1-phenyl-            | 88 J              |  |  |  |
| phenol, 2-ethyl-               | 24 J              |  |  |  |
| unknown (phthalate)            | 23 J              |  |  |  |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- J Indicates an estimated value less than the detection limit.

Job Number: ITEK 51037

#### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 5B-TF Lab Sample ID: SS8695

| Compound                   | <u>Concentration</u> |    | Compound                    | Concentr | ation |
|----------------------------|----------------------|----|-----------------------------|----------|-------|
| acenaphthene               | 20                   | U  | bis(2-chloroisopropyl)ether | 20       | U     |
| acenaphthylene             | 20                   | U  | bis(2-ethylhexyl)phthalate  | 5,000    | D     |
| anthracene                 | 20                   | U  | 4-bromophenyl phenyl ether  | 20       | U     |
| benzidine                  | 100                  | Ŭ. | 2-chloronaphthalene         | 20       | U     |
| benzo(a)anthracene         | 20                   | Ü: | 4-chlorophenyl phenyl ether | 20       | U     |
| benzo(b)fluoranthene       | 20                   | Ü  | chrysene                    | 20       | U     |
| benzo(k)fluoranthene       | 20                   | U, | dibenz(a,h)anthracene       | 20       | U     |
| benzo(a)pyrene             | 20                   | U: | di-n-butylphthalate         | 20       | U     |
| benzo(g,h,i)perylene       | 20                   | U  | 1,2-dichlorobenzene         | 20       | U     |
| butylbenzylphthalate       | 20                   | Ŭ. | 1,3-dichlorobenzene         | 20       | U     |
| bis(2-chloroethoxy)methane | 20                   | U  | 1,4-dichlorobenzene         | 20       | Ü     |
| bis(2-chloroethyl)ether    | 20                   | U  |                             |          |       |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction:

04/03/92

Date of Analysis:

04/20/92

J - Indicates an estimated value less than the detection limit.

D - Compound analyzed at a secondary dilution factor.

Job Number: ITEK 51037

#### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 5B-TF Lab Sample ID: SS8695

| Compound                  | Concentration | Compound                   | Concentration |  |  |
|---------------------------|---------------|----------------------------|---------------|--|--|
| 3,3'dichlorobenzidine     | <b>40</b> U   | hexachloroethane           | 20 U          |  |  |
| diethylphthalate          | 20 U          | indeno(1,2,3-cd)pyrene     | <b>20</b> Ú   |  |  |
| dimethyl phthalate        | 20 U          | isophorone                 | 20 U          |  |  |
| 2,4-dinitrotoluene        | 20 U          | naphthalene                | 20 U          |  |  |
| 2,6-dinitrotoluene        | 20 U          | nitrobenzene               | 20 U          |  |  |
| di-n-octyl phthalate      | 20 U          | n-nitroso-di-n-propylamine | 20 U          |  |  |
| 1,2-diphenylhydrazine(1)  | 20 U          | n-nitrosodimethylamine     | 20 U          |  |  |
| fluoranthene              | 20 U          | n-nitrosodiphenylamine(2)  | <b>20</b> U   |  |  |
| fluorene                  | 20 U          | phenanthrene               | <b>20</b> U   |  |  |
| hexachlorobenzene         | .20 U         | pyrene                     | 20 U          |  |  |
| hexachlorobutadiene       | 20 U          | 1,2,4-trichlorobenzene     | 20 U          |  |  |
| hexachlorocyclopentadiene | 20 U          | i ·                        |               |  |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

- Indicates an estimated value less than the detection limit.

Screened for as azobenzene

(2) - Detected as diphenylamine

Date of Extraction: Date of Analysis:

04/03/92

04/20/92

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

Client Project ID: LE Carpenter

Job Number: ITEK 51037

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 5B-TF Lab Sample ID: SS8695

| Tentative Identification (1)   | Concentration (2) |   |  |  |
|--------------------------------|-------------------|---|--|--|
| benzene, (1-methylethyl)-      | 41                | J |  |  |
| benzenemethanol, .alphamethyl- | 43                | J |  |  |
| ethanone, 1-phenyl-            | 66                | J |  |  |
| unknown (phthalate)            | 27                | J |  |  |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- J Indicates an estimated value less than the detection limit.

Job Number: ITEK 51037

#### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in  $\mu g/liter$  (ppb)

Sample Matrix: Water

Client Sample ID: 5C-TF Lab Sample ID: SS8696

| Compound                   | Concentration |    | Compound                    | Concentration |   |  |
|----------------------------|---------------|----|-----------------------------|---------------|---|--|
| acenaphthene               | 20            | U  | bis(2-chloroisopropyl)ether | 20            | U |  |
| acenaphthylene             | 20            | U  | bis(2-ethylhexyl)phthalate  | 9,200         | D |  |
| anthracene                 | 20            | U  | 4-bromophenyl phenyl ether  | 20            | U |  |
| benzidine                  | 100           | U  | 2-chloronaphthalene         | 20            | U |  |
| benzo(a)anthracene         | 20            | U  | 4-chlorophenyl phenyl ether | 20            | U |  |
| benzo(b)fluoranthene       | 20            | U, | chrysene                    | 20            | U |  |
| benzo(k)fluoranthene       | 20            | U  | dibenz(a,h)anthracene       | 20            | U |  |
| benzo(a)pyrene             | 20            | U  | di-n-butylphthalate         | 20            | U |  |
| benzo(g,h,i)perylene       | 20            | U  | 1,2-dichlorobenzene         | 20            | U |  |
| butylbenzylphthalate       | 20            | U  | 1,3-dichlorobenzene         | ,20           | U |  |
| bis(2-chloroethoxy)methane | 20            | U  | 1,4-dichlorobenzene         | 20            | Ů |  |
| bis(2-chloroethyl)ether    | 20            | U  |                             |               |   |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 04/03/92 Date of Analysis: 04/20/92

J - Indicates an estimated value less than the detection limit.

D - Compound analyzed at a secondary dilution factor.

Job Number: ITEK 51037

### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 5C-TF Lab Sample ID: SS8696

| • | Compound                  | Concentration |          | Compound                   | Concentration |   |  |
|---|---------------------------|---------------|----------|----------------------------|---------------|---|--|
|   | 3,3'dichlorobenzidine     | 40            | Ü        | hexachloroethane           | 20            | Ų |  |
|   | diethylphthalate          | 20            | <b>U</b> | indeno(1,2,3-cd)pyrene     | 20            | U |  |
|   | dimethyl phthalate        | 20            | Ü        | isophorone                 | 20            | U |  |
|   | 2,4-dinitrotoluene        | 20            | U        | naphthalene                | 20            | U |  |
|   | 2,6-dinitrotoluene        | 20            | U        | nitrobenzene               | 20            | U |  |
|   | di-n-octyl phthalate      | 6             | J        | n-nitroso-di-n-propylamine | 20            | U |  |
|   | 1,2-diphenylhydrazine(1)  | 20            | Ü        | n-nitrosodimethylamine     | 20            | U |  |
|   | fluoranthene              | 20            | Ü        | n-nitrosodiphenylamine(2)  | 20            | Ü |  |
|   | fluorene                  | 20            | U        | phenanthrene               | 20            | U |  |
|   | hexachlorobenzene         | 20            | U        | pyrene                     | 20            | U |  |
|   | hexachlorobutadiene       | 20            | U        | 1,2,4-trichlorobenzene     | 20            | U |  |
|   | hexachlorocyclopentadiene | 20            | U        |                            |               |   |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

(1) - Screened for as azobenzene

(2) - Detected as diphenylamine

Date of Extraction: 04/03/92 Date of Analysis: 04/20/92

Job Number: ITEK 51037

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in  $\mu g/\text{liter}$  (ppb)

Sample Matrix: Water

Client Sample ID: 5C-TF Lab Sample ID: SS8696

| Tentative Identification (1)                 |   | Concentration (2) |   |  |
|--|---|-------------------|---|--|
| benzene, (1-methylethyl)-                    | + | 61                | j |  |
| benzenemethanol, .alphamethyl-               |   | 32                | J |  |
| ethanone, 1-phenyl-                          | • | 99                | J |  |
| phenol, 2-ethyl-                             |   | 19                | J |  |
| phosphoric acid, 2-ethylhexyl diphenyl ester |   | 16                | J |  |
| unknown (phthalate)                          |   | 61                | J |  |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- J Indicates an estimated value less than the detection limit.

KNOXVILLE, TN

Job Number: ITEK 51037

#### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank Lab Sample ID: BL0203

| Compound                   | Concentration | Compound                    | Concentration |
|----------------------------|---------------|-----------------------------|---------------|
| acenaphthene               | 10 U          | bis(2-chloroisopropyl)ether | 10 U          |
| acenaphthylene             | 10 U          | bis(2-ethylhexyl)phthalate  | 10 U          |
| anthracene                 | 10 U          | 4-bromophenyl phenyl ether  | 10 U          |
| benzidine                  | 50 U          | 2-chloronaphthalene         | 10 U          |
| benzo(a)anthracene         | 10 U          | 4-chlorophenyl phenyl ether | 10 U          |
| benzo(b)fluoranthene       | 10 U          | chrysene                    | 10 U          |
| benzo(k)fluoranthene       | 10 U          | dibenz(a,h)anthracene       | 10 U          |
| benzo(a)pyrene             | 10 U          | di-n-butylphthalate         | 10 U          |
| benzo(g,h,i)perylene       | 10 U          | 1,2-dichlorobenzene         | 10 U          |
| butylbenzylphthalate       | 10 U          | 1,3-dichlorobenzene         | 10 U          |
| bis(2-chloroethoxy)methane | 10 U          | 1,4-dichlorobenzene         | 10 U          |
| bis(2-chloroethyl)ether    | 10 U          |                             |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 04/03/92 Date of Analysis: 04/06/92

This method blank applies to the following samples: 1A-TF, 1B-TF, 1C-TF, 2A-TF, 2A-TF DL, 2B-TF, 2B-TF DL, 2C-TF, 2C-TF DL, 3A-TF, 3B-TF, 3C-TF, 4A-TF, 4A-TF DL, 4B-TF, 4B-TF DL, 4C-TF, 4C-TF DL, 5A-TF, 5A-TF DL, 5B-TF, 5B-TF DL, 5C-TF and 5C-TF DL.

DL = Dilution

J - Indicates an estimated value less than the detection limit.

Job Number: ITEK 51037

### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank Lab Sample ID: BL0203

| Compound                  | Concentration | Compound                   | Concentration |
|---------------------------|---------------|----------------------------|---------------|
| 3,3'dichlorobenzidine     | 20 U          | hexachloroethane           | 10 U          |
| diethylphthalate          | 10 U          | indeno(1,2,3-cd)pyrene     | 10 U          |
| dimethyl phthalate        | 10 U          | isophorone                 | 10 U          |
| 2,4-dinitrotoluene        | 10 U          | naphthalene                | 10 U          |
| 2,6-dinitrotoluene        | 10 U          | nitrobenzene               | 10 U          |
| di-n-octyl phthalate      | 10 U          | n-nitroso-di-n-propylamine | 10 U          |
| 1,2-diphenylhydrazine(1)  | 10 U          | n-nitrosodimethylamine     | 10 U          |
| fluoranthene              | 10 U          | n-nitrosodiphenylamine(2)  | 10 U          |
| fluorene                  | 10 U          | phenanthrene               | 10 U          |
| hexachlorobenzene         | 10 U          | pyrene                     | 10 U          |
| hexachlorobutadiene       | 10 U          | 1,2,4-trichlorobenzene     | 10 U          |
| hexachlorocyclopentadiene | 10 U          |                            |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 04/03/92 Date of Analysis: 04/06/92

J - Indicates an estimated value less than the detection limit.

<sup>(1) -</sup> Screened for as azobenzene

<sup>(2) -</sup> Detected as diphenylamine

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

Client Project ID: LE Carpenter

Job Number: ITEK 51037

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in  $\mu g/liter$  (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank Lab Sample ID: BL0203

Tentative Identification (1)

2-pentanone, 4-hydroxy-4-methylunknown

Concentration (2)

11 JA

12 JA

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- A Suspected aldol product.
- J Indicates an estimated value less than the detection limit.

Job Number: ITEK 51037

#### WATER SURROGATE PERCENT RECOVERY SUMMARY

#### **SEMI-VOLATILE**

|                  |                                    | (C)                                 | A Section of the section of     |                     |                                   | 2,4,6-                            |
|------------------|------------------------------------|-------------------------------------|---------------------------------|---------------------|-----------------------------------|-----------------------------------|
| Client Sample ID | Nitro-<br>Benzene-D5<br>(35-114%)* | 2-Fluoro-<br>Biphenyl<br>(43-116%)* | Terphenyl-<br>D14<br>(33-141%)* | Phenol-D5 (10-94%)* | 2-Fluoro-<br>Phenol<br>(21-100%)* | Tribromo-<br>Phenol<br>(10-123%)* |
| 1A-TF            | 80                                 | 83                                  | 64                              | NA                  | NA                                | NA                                |
| 1B-TF            | 77                                 | 82                                  | 64                              | NA                  | NA                                | NA                                |
| 1C-TF            | 78                                 | 85                                  | 72                              | NA                  | NA                                | NA <sub>.</sub>                   |
| 2A-TF            | 71                                 | 69                                  | 32 **                           | NA                  | NA ·                              | NA                                |
| 2A-TF DL         | 70                                 | 65                                  | 30 **                           | NA                  | NA                                | NA                                |
| 2B-TF            | <b>79</b>                          | 71                                  | 34                              | NA                  | NA                                | NA                                |
| 2B-TF DL         | 77                                 | 65                                  | 31 **                           | NA                  | NA.                               | NA                                |
| 2C-TF            | 86                                 | 79                                  | 40                              | NA                  | NA                                | ŅA                                |
| 2C-TF DL         | 75                                 | 73                                  | 38                              | NA                  | NA                                | NA                                |
| 3A-TF            | 85                                 | 85                                  | 65                              | NA                  | NA                                | NA                                |
| 3B-TF            | 82                                 | 82                                  | 64                              | NA                  | NA                                | NA                                |
| 3C-TF            | 78                                 | 78                                  | 64                              | NA                  | NA                                | NA                                |
| 4A-TF            | 82                                 | 85                                  | 59                              | NA                  | NA.                               | NA.                               |
| 4A-TF DL         | D                                  | D                                   | D                               | NA                  | NA                                | NA                                |
| 4B-TF            | 85                                 | 88                                  | 55                              | ŇA                  | NA                                | NA                                |
| 4B-TF DL         | D .                                | D                                   | D                               | NA                  | NA                                | NA                                |
| 4C-TF            | 78                                 | 84                                  | 63                              | NA                  | NA                                | NA                                |
| 4C-TF DL         | D                                  | D                                   | D.                              | NA                  | NA                                | NA                                |

<sup>\*</sup>Values in parenthesis represent QC limits.

D - Surrogates diluted out

NA - Not applicable

DL - Dilution

<sup>\*\*</sup>Values are outside of QC limits.

Job Number: ITEK 51037

#### WATER SURROGATE PERCENT RECOVERY SUMMARY

#### **SEMI-VOLATILE**

| Client Sample ID | Nitro-<br>Benzene-D5<br>(35-114%)* | 2-Fluoro-<br>Biphenyl<br>(43-116%)* | Terphenyl-<br>D14<br>(33-141%)* | Phenol-D5<br>(10-94%)* | 2-Fluoro-<br>Phenol<br>(21-100%)* | 2,4,6-<br>Tribromo-<br>Phenol<br>(10-123%)* |
|------------------|------------------------------------|-------------------------------------|---------------------------------|------------------------|-----------------------------------|---|
| 5A-TF            | 79                                 | 85                                  | 49                              | NA                     | NA                                | NA  |
| 5A-TF DL         | D                                  | <b>D</b> )                          | D                               | NA                     | NA                                | NA  |
| 5B-TF            | 77                                 | 84                                  | 44                              | NA                     | NA                                | NA .  |
| 5B-TF DL         | D                                  | D                                   | D                               | NA                     | NA                                | NA  |
| 5C-TF            | 75                                 | 85                                  | 54                              | NA                     | NA                                | NA  |
| 5C-TF DL         | D                                  | D                                   | D                               | NA                     | NA                                | NA  |
| Method Blank     | 94                                 | .95                                 | <del>9</del> 6                  | NA                     | NA                                | NA  |

\*Values in parenthesis represent QC limits.

D - Surrogates diluted out

NA - Not applicable

DL - Dilution

# Appendix D

Geotechnical Data



### ANALYTICAL **SERVICES**

### CERTIFICATE OF ANALYSIS

Michael Krstich IT Corporation 312 Directors Drive Knoxville, Tn. 37923

May 14, 1992

ETDC Project Number: 483500.004.01

P.O. Number: 483048

This is the Certificate of Analysis for the following samples:

Client Project ID: Date Received by Lab:

L.E. Carpenter April 8, 1992

Number of Samples:

Six (6)

Sample Type:

Soil

#### I. Introduction/Case Narrative

Six soil samples were received by IT/ETDC on April 8, 1992 for analysis of grainsize distribution, moisture content, unit weight density, organic content, atterberg limits, specific gravity, permeability and porosity.

Please see Appendix A, the sample number cross reference list; Appendix B, the analysis results; and Appendix C, the Chain of Custody and Request for Analysis records.

Reviewed and Approved:

Beverly E. Heamon

Manager, Geotechnical Services

Page 2 of 16 Michael Krstich IT Corporation May 14, 1992 Client Project I IT ANALYTICAL SERVICES
304 DIRECTORS DRIVE
KNOXVILLE, TN
(615) 690-3211

Client Project ID: L.E. Carpenter ETDC Project No.: 483500.004.01

#### II. Analytical Results/Methodology

REFERENCES: Army Corps of Engineers laboratory Soils Testing, EM1110-2-1906; EPA SW-846, Test Methods for Evaluating Solid Waste. Annual Book of ASTM Standards, Section 4, Construction, Volume 04.08, Soil and Rock; Dimension Stone; Geosynthetics. Volume 4.02, Concrete and Aggregates.

Unit Weight Density & Porosity
Permeabiltiy
Grainsize Distribution
Moisture Content
Atterberg Limits
Ash & Organic Content
Specific Gravity

EM1110-2-1906 8W846, Method 9100 ASTM D 422 ASTM D 2216 ASTM D 4318 ASTM D 2974 ASTM D 854/ASTM C 127

#### III. Quality Control

Quality control checks such as duplicates and spikes (QC samples), are not normally applicable to geotechnical testing. This is due to the inability of obtaining samples with known characteristics, the heterogenous nature of the samples, and Quality Control procedures built-in to the analytical method.

QC measures to ensure accuracy and precision of test results include the following:

- 100% verification on all numerical results all raw data entries, transcriptions and calculations entered by lab technicians are checked, recalculated and verified. Most data calculations are performed by computer programs.
- Data validation through test reasonableness summaries of all test results for individual reports are reviewed to determine the overall reasonableness of data and to determine the presence of any data that may be considered outliers.
- Quality control procedures are built into most standardized geotechnical procedures. For example, many analyses routinely call for a re-analysis, specifying an acceptance criteria.
- Routine instrument calibration all instruments, gauges and equipment used in testing is calibrated on a routine basis. All instrument calibration follows ASTM or manufacturer guidelines.

Page 3 of 16 Michael Krstich IT Corporation May 14, 1992 Client Project I IT ANALYTICAL SERVICES 304 DIRECTORS DRIVE KNOXVILLE, TN (615) 690-3211

Client Project ID: L.E. Carpenter ETDC Project No.: 483500.004.01

- Maintenance of all past calibration records records and certification documents of all instruments, gauges and equipment are updated routinely and maintained in the Quality Control Coordinators Quality/Operations files.
- Use of trained personnel for conducting tests all technicians are trained in the application of standard laboratory procedures for geotechnical analyses as well as the quality assurance measures implemented by IT.

#### IV. Data Qualification

Michael Krstich of IT/Knoxville requested geotechnical test parameters on six of the thirteen samples received by IT/ETDC. Samples were sent to IT/ETDC as a group, so they could be returned together to their site of origin after geotechnical testing was completed. Geotechnical testing was requested and completed on client sample numbers ST-2A/24-48, ST-2A/48-72, ST-2B/24-48, ST-2B/72-96, ST-3A/0-24 and ST-3B/24-48 (ETDC numbers: 1993-1995 and 1997-1999). Client sample number ST-2B/48-72 (ETDC-1996) lacked sufficient sample quantity to complete all of the originally requested geotechnical test parameters. With the knowledge and approval of Michael Krstich, client sample number ST-3B/24-48 (ETDC-1999) was substituted in place of client sample number ST-2B/48-72 (ETDC-1996) for geotechnical analysis testing.

Average specific gravity values are calculated based on an average of the bulk specific gravity\* and the specific gravity of material passing the #10 sieve. Average specific gravity values given on most data results mean that the sample was composed of particles which were smaller than the #4 sieve and a signicant amount of material which was greater than the #4 sieve. Client sample number ST-3B/24-48 (ETDC-1999) was the only sample which did not contain a signicant amount of material greater than the #4 sieve. The specific gravity value given on data results for ETDC-1999 represents the specific gravity of material passing the #10 sieve. The average specific gravity calculation is referenced from ASTM D 854 and ASTM C 127.

Fine sieve and hydrometer results occasionally overlap due to organic debris, soluble salts or other contaminants contained in the sample. Data points are plotted as calculated. No attempt has been made to curve-fit the grainsize data points.

\*Bulk specific gravity is defined by the specific gravity of the material greater than the number 4 sieve (4.75 mm).

Page 4 of 16
Michael Krstich
IT Corporation
May 14, 1992
Client Project ID: L.E. Carpenter

IT ANALYTICAL SERVICES 304 DIRECTORS DRIVE KNOXVILLE, TN (615) 690-3211

ETDC Project No.: 483500.004.01

#### CROSS-REFERENCE LIST

| ETDC SAMPLE NO. | CLIENT SAMPLE NO. |
|-----------------|-------------------|
|                 |                   |
| ETDC-1993       | ST-2A/24-48       |
|                 |                   |
|                 |                   |
| ETDC-1997       | <br>ST-2B/72-96   |
| ETDC-1998       | <br>ST-3A/0-24    |
|                 |                   |

Page 5 of 16 Michael Krstich IT Corporation May 14, 1992 IT ANALYTICAL SERVICES 304 DIRECTORS DRIVE KNOXVILLE, TN (615) 690-3211

Client Project ID: L.E. Carpenter ETDC Project No.: 483500.004.01

#### SAMPLE ANALYSIS RESULTS

PROJECT NAME: L.E. CARPENTER USCS SYMBOL: ML WATER CONTENT, %: PROJECT NO.: 483500.004.01 6.5 CUST. SAMPLE NO.: ST-2A/24-48 LIQUID LIMIT: 26.0 ETDC SAMPLE NO.: ETDC-1993 PLASTICITY INDEX: 2.0 AVG. SPECIFIC GRAVITY: 2.8201 (MEASURED) PERMEABILITY: 5.1 E-5 cm/s UNIT WEIGHT DENSITY: 122.6 pcf ORGANIC CONTENT, %: 5.3 ASH CONTENT, %: 94.7 DRY DENSITY: 115.1 pcf POROSITY: 34.6

=====SIEVE ANALYSIS==========

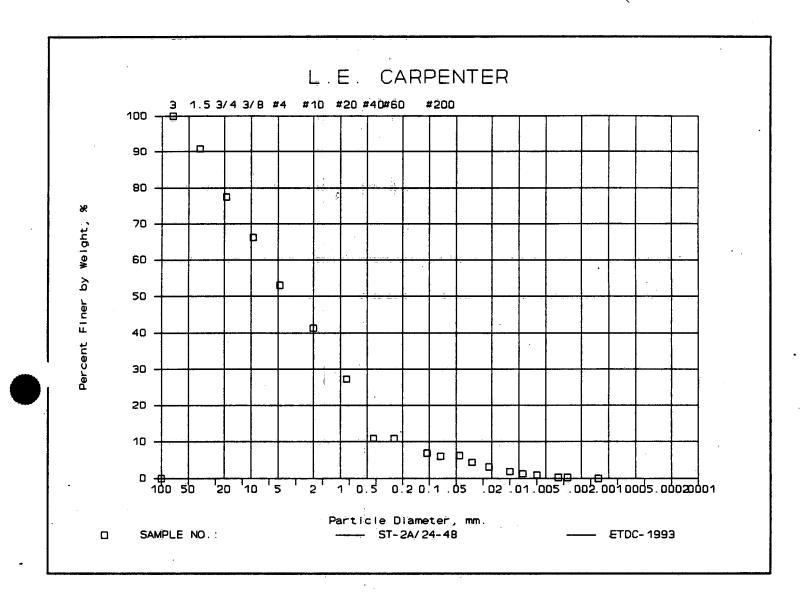
| SIEVE NO. | DIAMETER (mm) | PERCENT FINER (%) |
|-----------|---------------|-------------------|
|           |               |                   |
| 3.0 in    | 75.000        | 100.0             |
| 1.5 in    | 37.500        | 90.8              |
| 0.75 in   | 19.000        | 77.5              |
| 0.375 in  | 9.500         | 66.2              |
| NO. 4     | 4.750         | 53.0              |
| NO. 10    | 2.000         | 41.4              |
| NO. 20    | 0.850         | 27.2              |
| NO. 40    | 0.425         | 10.9              |
| NO. 60    | 0.250         | 10.9              |
| NO. 140   | 0.106         | 6.9               |
| NO. 200   | 0.075         | 6.0               |

-----HYDROMETER ANALYSIS----

| PERCENT FINER |
|---------------|
| (%)           |
|               |
| 6.2           |
| 4.3           |
| 3.1           |
| 1.8           |
| 1.2           |
| 0.9           |
| 0.3           |
| 0.3           |
| 0.0           |
|               |

Page 6 of 16 Michael Krstich IT Corporation May 14, 1992 IT ANALYTICAL SERVICES 304 DIRECTORS DRIVE KNOXVILLE, TN (615) 690-3211

Client Project ID: L.E. Carpenter ETDC Project No.: 483500.004.01



Page 7 of 16 Michael Krstich IT Corporation May 14, 1992 IT ANALYTICAL SERVICES 304 DIRECTORS DRIVE KNOXVILLE, TN (615) 690-3211

May 14, 1992 Client Project ID: L.E. Carpenter ETDC Project No.: 483500.004.01

#### SAMPLE ANALYSIS RESULTS

| PROJECT NAME:          | L.E. CARPENTER    | USCS SYMBOL:         | *NP          |
|------------------------|-------------------|----------------------|--------------|
| PROJECT NO.:           | 483500.004.01     | WATER CONTENT, %:    | 15.1         |
| CUST. SAMPLE NO.:      | ST-2A/48-72       | LIQUID LIMIT:        | 30.0         |
| ETDC SAMPLE NO.:       | ETDC-1994         | PLASTICITY INDEX:    | *NP          |
| AVG. SPECIFIC GRAVITY: | 2.5610 (MEASURED) | PERMEABILITY:        | 1.8 E-5 cm/s |
| ORGANIC CONTENT, %:    | 8.0               | UNIT WEIGHT DENSITY: | 60.6 pcf     |
| ASH CONTENT, %:        | 92.0              | DRY DENSITY:         | 52.6 pcf     |
| POROSITY:              | 67.1              |                      | , -          |

#### ----SIEVE ANALYSIS-----

| SIEVE NO. | DIAMETER (mm) | PERCENT FINER (%) |
|-----------|---------------|-------------------|
|           |               |                   |
| 3.0 in    | 75.000        | 100.0             |
| 1.5 in    | 37.500 °      | 100.0             |
| 0.75 in   | 19.000        | 72.4              |
| 0.375 in  | 9.500         | 71.7              |
| NO. 4     | 4.750         | 64.2              |
| NO. 10    | 2.000         | 60.0              |
| NO. 20    | 0.850         | 51.3              |
| NO. 40    | 0.425         | 34.4              |
| NO. 60    | 0.250         | 34.4              |
| NO. 140   | 0.106         | 25.0              |
| NO. 200   | 0.075         | 23.0              |

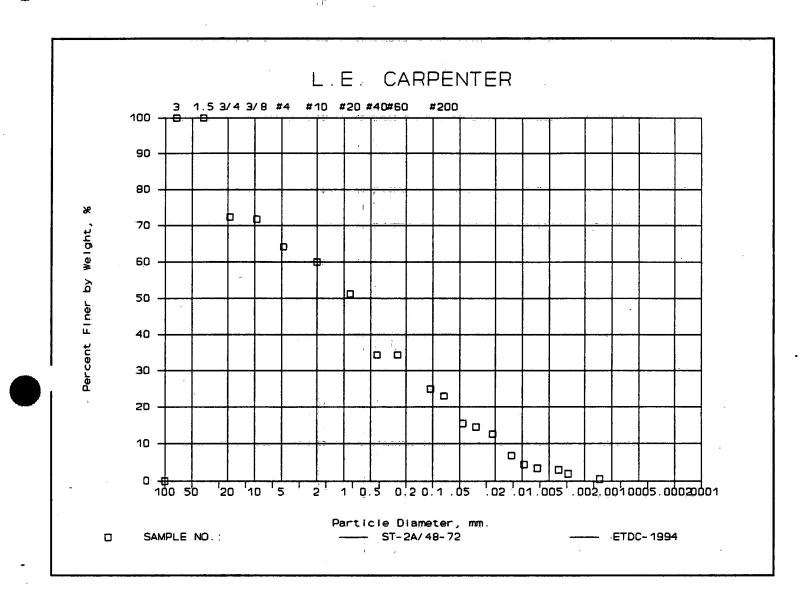
#### ----HYDROMETER ANALYSIS-

| DIAMETER (mm) | PERCENT FINER (%) |
|---------------|-------------------|
|               | *******           |
| 0.0460        | 15.5              |
| 0.0328        | 14.5              |
| 0.0213        | 12.6              |
| 0.0131        | 6.8               |
| 0.0095        | 4.4               |
| 0.0067        | 3.4               |
| 0.0039        | 2.9               |
| 0.0031        | 1.9               |
| 0.0014        | 0.5               |

*NPLASTIC* 

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Client Project ID: L.E. Carpenter ETDC Project No.: 483500.004.01



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May 14, 1992
Client Project ID: L.E. Carpenter ETDC Project No.: 483500.004.01

#### SAMPLE ANALYSIS RESULTS

| PROJECT NAME:          | L.E. CARPENTER    | USCS SYMBOL:         | *NP          |
|------------------------|-------------------|----------------------|--------------|
| PROJECT NO.:           | 483500.004.01     | WATER CONTENT, %:    | 6.6          |
| CUST. SAMPLE NO.:      | ST-2B/24-48       | LIQUID LIMIT:        | 24.0         |
| ETDC SAMPLE NO.:       | ETDC-1995         | PLASTICITY INDEX:    | *NP          |
| AVG. SPECIFIC GRAVITY: | 2.9551 (MEASURED) | PERMEABILITY:        | 1.1 E-4 cm/s |
| ORGANIC CONTENT, %:    | 4.2               | UNIT WEIGHT DENSITY: | 99.7 pcf     |
| ASH CONTENT, %:        | 95.8              | DRY DENSITY:         | 93.5 pcf     |
| POROSITY:              | 49.3              | ,                    | -            |

#### ====SIEVE ANALYSIS===

| SIEVE NO. | DIAMETER<br>(mm) | PERCENT FINER (%) |
|-----------|------------------|-------------------|
|           |                  |                   |
| 3.0 in    | 75.000           | 100.0             |
| 1.5 in    | 37.500           | 100.0             |
| 0.75 in   | 19.000           | 86.5              |
| 0.375 in  | 9.500            | 70.8              |
| NO. 4     | 4.750            | 61.2              |
| NO. 10    | 2.000            | 40.6              |
| NO. 20    | 0.850            | 29.5              |
| NO. 40    | 0.425            | 15.6              |
| NO. 60    | 0.250            | 15.5              |
| NO. 140   | 0.106            | 10.1              |
| NO. 200   | 0.075            | 8.8               |

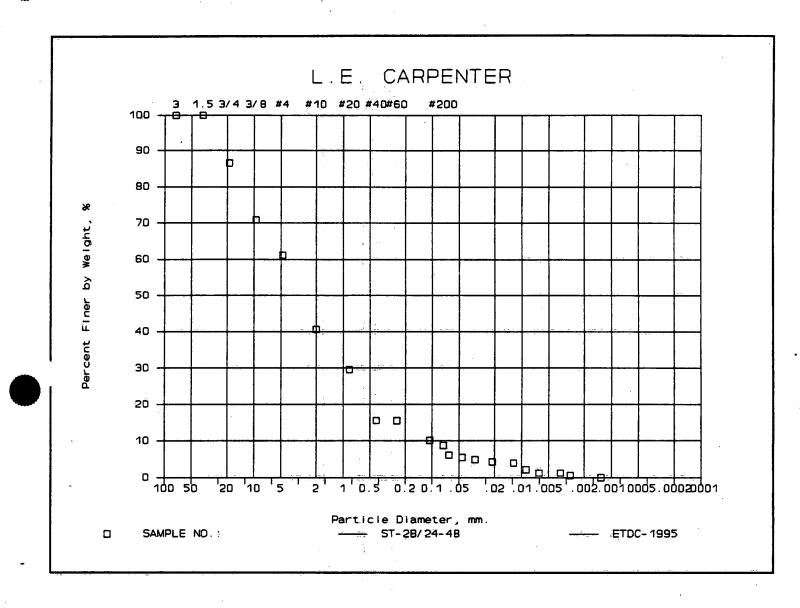
#### ======HYDROMETER ANALYSIS====

| DIAMETER (mm) | PERCENT FINER (%) |
|---------------|-------------------|
|               |                   |
| 0.0647        | 6.1               |
| 0.0461        | 5.5               |
| 0.0330        | 4.9               |
| 0.0211        | 4.3               |
| 0.0122        | 4.0               |
| 0.0089        | 2.1               |
| 0.0064        | 1.2               |
| 0.0037        | 1.2               |
| 0.0029        | 0.6               |
| 0.0013        | 0.0               |

**UNPLASTIC** 

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Client Project ID: L.E. Carpenter ETDC Project No.: 483500.004.01



IT ANALYTICAL SERVICES 304 DIRECTORS DRIVE KNOXVILLE, TN (615) 690-3211

May 14, 1992 Client Project ID: L.E. Carpenter ETDC Project No.: 483500.004.01

#### SAMPLE ANALYSIS RESULTS

| PROJECT NAME:          | L.E. CARPENTER    | USCS SYMBOL:         | ML           |
|------------------------|-------------------|----------------------|--------------|
| PROJECT NO.:           | 483500.004.01     | WATER CONTENT, %:    | 5.6          |
| CUST. SAMPLE NO.:      | ST-2B/72-96       | LIQUID LIMIT:        | 23.0         |
| ETDC SAMPLE NO.:       | ETDC-1997         | PLASTICITY INDEX:    | 1.0          |
| AVG. SPECIFIC GRAVITY: | 2.8363 (MEASURED) | PERMEABILITY:        | 1.4 E-4 cm/s |
| ORGANIC CONTENT, %:    | 3.8               | UNIT WEIGHT DENSITY: | 106.0 pcf    |
| ASH CONTENT, %:        | 96.2              | DRY DENSITY:         | 100.3 pcf    |
| POROSITY:              | 43.3              |                      | - · ·        |

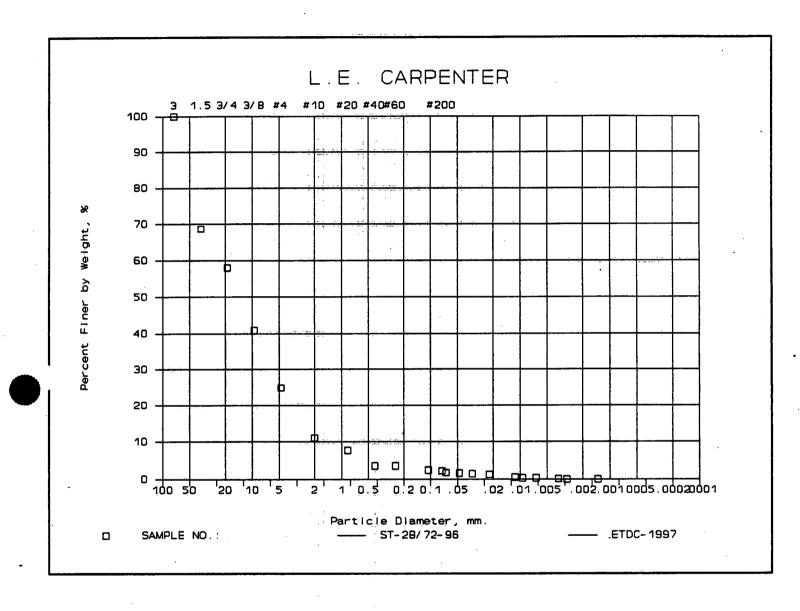
------SIEVE ANALYSIS------

| SIEVE NO. | DIAMETER (mm) | PERCENT FINER (%) |
|-----------|---------------|-------------------|
|           |               |                   |
| 3.0 in    | 75.000        | 100.0             |
| 1.5 in    | 37.500        | 68.7              |
| 0.75 in   | 19.000        | 58.0              |
| 0.375 in  | 9.500         | 41.0              |
| NO. 4     | 4.750         | 24.8              |
| NO. 10    | 2.000         | 11.0              |
| NO. 20    | 0.850         | 7.6               |
| NO. 40    | 0.425         | 3.6               |
| NO. 60    | 0.250         | 3.6               |
| NO. 140   | 0.106         | 2.4               |
| NO. 200   | 0.075         | 2.2               |

| DIAMETER (mm) | PERCENT FINER (%) |
|---------------|-------------------|
|               |                   |
| 0.0675        | 1.7               |
| 0.0479        | 1.6               |
| 0.0342        | 1.5               |
| 0.0220        | 1.2               |
| 0.0114        | 0.5               |
| 0.0094        | 0.3               |
| 0.0066        | 0.3               |
| 0.0037        | 0.1               |
| 0.0030        | 0.0               |
| 0.0014        | 0.0               |

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Client Project ID: L.E. Carpenter ETDC Project No.: 483500.004.01



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May 14, 1992
Client Project ID: L.E. Carpenter ETDC Project No.: 483500.004.01

#### SAMPLE ANALYSIS RESULTS

| PROJECT NAME:          | L.E. CARPENTER    | USCS SYMBOL:         | CL           |
|------------------------|-------------------|----------------------|--------------|
| PROJECT NO.:           | 483500.004.01     | WATER CONTENT, %:    | 10.2         |
| CUST. SAMPLE NO.:      | ST-3A/0-24        | LIQUID LIMIT:        | 20.0         |
| ETDC SAMPLE NO.:       | ETDC-1998         | PLASTICITY INDEX:    | 3.0          |
| AVG. SPECIFIC GRAVITY: | 2.6320 (MEASURED) | PERMEABILITY:        | 3.3 E-5 cm/s |
| ORGANIC CONTENT, %:    | 3.4               | UNIT WEIGHT DENSITY: | 111.3 pcf    |
| ASH CONTENT, %:        | 96.6              | DRY DENSITY:         | 101.0 pcf    |
| POROSITY:              | 38.5              |                      |              |

----SIEVE ANALYSIS-----

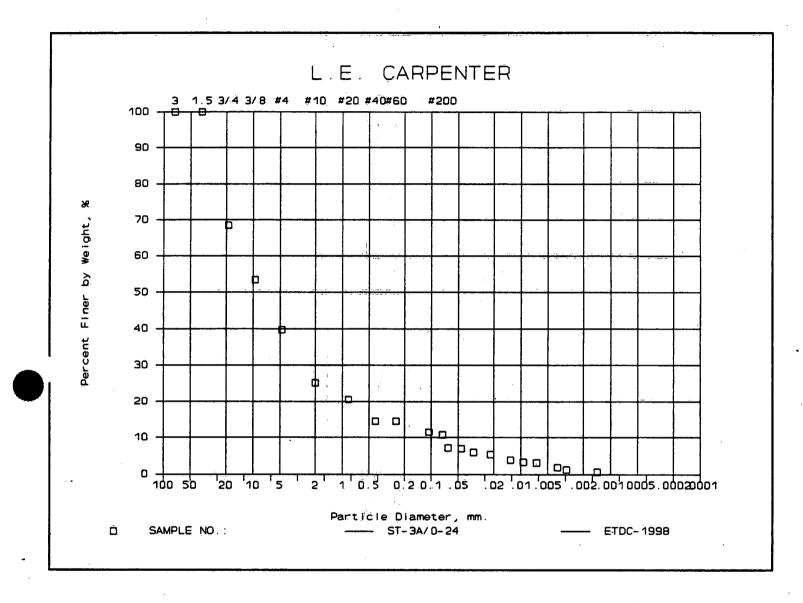
| SIEVE NO. | DIAMETER (mm) | PERCENT FINER (%) |
|-----------|---------------|-------------------|
|           | <del></del>   | 100.0             |
| 3.0 in    | 75.000        | 100.0             |
| 1.5 in    | 37.500        | 100.0             |
| 0.75 in   | 19.000        | 68.5              |
| 0.375 in  | 9.500         | 53.5              |
| NO. 4     | 4.750         | 39.7              |
| NO. 10    | 2.000         | 25.1              |
| NO. 20    | 0.850         | 20.6              |
| NO. 40    | 0.425         | 14.5              |
| NO. 60    | 0.250         | 14.5              |
| NO. 140   | 0.106         | 11.4              |
| NO. 200   | 0.075         | 10.8              |

#### 

| DIAMETER (mm) | PERCENT FINER (%) |
|---------------|-------------------|
| 0.0648        | 7.1               |
| 0.0461        | 6.9               |
| 0.0336        | 5.9               |
| 0.0216        | 5.3               |
| 0.0129        | 3.9               |
| 0.0093        | 3.3               |
| 0.0066        | 3.1               |
| 0.0039        | 1.8               |
| 0.0031        | 1.2               |
| 0.0014        | 0.6               |

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May 14, 1992 Client Project ID: L.E. Carpenter ETDC Project No.: 483500.004.01



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IT ANALYTICAL SERVICES 304 DIRECTORS DRIVE KNOXVILLE, TN (615) 690-3211

\*NP

40.2

30.0

Client Project ID: L.E. Carpenter ETDC Project No.: 483500.004.01

#### SAMPLE ANALYSIS RESULTS

PROJECT NAME: PROJECT NO.: CUST. SAMPLE NO.: ETDC SAMPLE NO.: SPECIFIC GRAVITY: ORGANIC CONTENT, %: ASH CONTENT, %: POROSITY:

L.E. CARPENTER 483500.004.01 ST-3B/24-48 ETDC-1999

2.5174 (MEASURED) 4.3 95.7

46.8

LIQUID LIMIT: PLASTICITY INDEX: PERMEABILITY:

USCS SYMBOL:

WATER CONTENT, %:

\*NP 5.6 E-6 cm/s

UNIT WEIGHT DENSITY: 117.2 pcf DRY DENSITY: 83.6 pcf

====SIEVE ANALYSIS=====

| SIEVE NO. | DIAMETER | PERCENT FINER |
|-----------|----------|---------------|
|           | (mm)     | (%)           |
|           |          |               |
| 3.0 in    | 75.000   | 100.0         |
| 1.5 in    | 37.500   | 100.0         |
| 0.75 in   | 19.000   | 100.0         |
| 0.375 in  | 9.500    | 100.0         |
| NO. 4     | 4.750    | 99.2          |
| NO. 10    | 2.000    | 97.6          |
| NO. 20    | 0.850    | 87.3          |
| NO. 40    | 0.425    | 75.7          |
| NO. 60    | 0.250    | 68.7          |
| NO. 140   | 0.106    | 58.2          |
| NO. 200   | 0.075    | 56.7          |

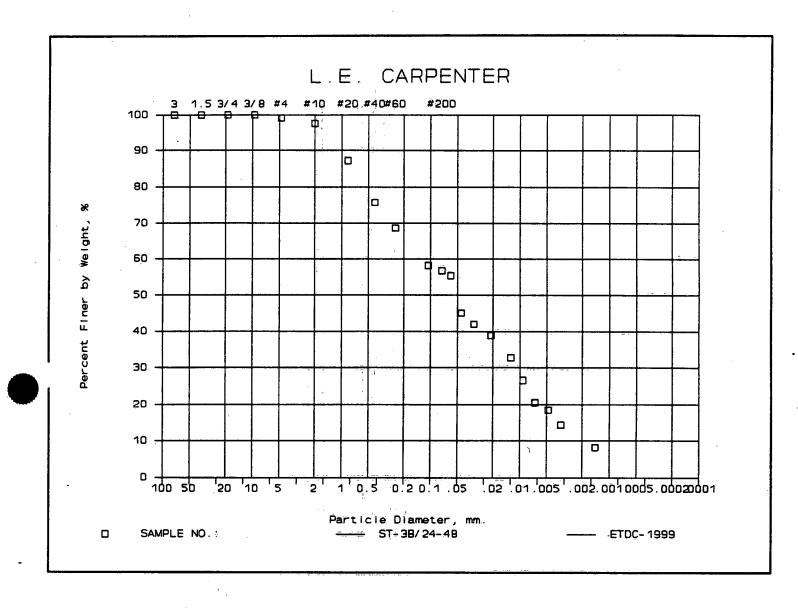
#### =HYDROMETER ANALYSIS====

| DIAMETER<br>(mm) | PERCENT FINER (%) |
|------------------|-------------------|
|                  |                   |
| 0.0594           | 55.3              |
| 0.0451           | 45.1              |
| 0.0326           | 42.0              |
| 0.0210           | 38.9              |
| 0.0126           | 32.8              |
| 0.0092           | 26.6              |
| 0.0067           | 20.5.             |
| 0.0048           | 18.4              |
| 0.0035           | 14.3              |
| 0.0014           | 8.2               |
|                  |                   |

ONPLASTIC

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Client Project ID: L.E. Carpenter ETDC Project No.: 483500.004.01



Job Number: ITAD 51191

#### TOTAL ORGANIC CARBON ANALYSIS

Results in mg/kg (ppm)

Sample Matrix: Soil

| Client Sample ID | Lab Sample ID | Result |   |
|------------------|---------------|--------|---|
| Method Blank     | P3763         | 1 U    | ļ |
| ST-2A/24-48      | K0494         | 18,000 |   |
| ST-2B/72-96      | K0495         | 3,500  |   |
| ST-2A/48-72      | K0496         | 2,300  |   |
| ST-2B/24-48      | K0497         | 3,300  |   |
| ST-2B/48-72      | K0498         | 8,800  |   |
| ST-3A/0-24       | K0499         | 4,100  |   |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 04/22/92

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

Client Project ID: LE Carpenter

Job Number: ITAD 51328

#### TOTAL ORGANIC CARBON ANALYSIS

Results in mg/kg (ppm)

Sample Matrix: Soil

 Client Sample ID
 Lab Sample ID
 Result

 Method Blank
 P3825
 1 U

 ST-3B/24-48
 K1791
 7,000

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 05/19/92

### Appendix E

Soils Analyses

Job Number: ITAD 51019

# VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS Results in μg/kg (ppb) Sample Matrix: Soil

Client Sample ID: 1045-052-1 Lab Sample ID: \$\$8495

|                          |         |        | ,                         |         |        |
|--------------------------|---------|--------|---------------------------|---------|--------|
| Compound                 | Concent | ration | Compound                  | Concent | ration |
| acrolein                 | 12,000  | U      | 1,1-dichloroethene        | 6,200   | U      |
| acrylonitrile            | 12,000  | Ü      | trans-1,2-dichloroethene  | 6,200   | U      |
| benzene                  | 6,200   | U      | 1,2-dichloropropane       | 6,200   | U      |
| bromodichloromethane     | 6,200   | U      | cis-1,3-dichloropropene   | 6,200   | U      |
| bromoform                | 6,200   | U      | trans-1,3-dichloropropene | 6,200   | U      |
| bromomethane             | 12,000  | U      | ethyl benzene             | 130,000 |        |
| carbon tetrachloride     | 6,200   | U      | methylene chloride        | 6,200   | U      |
| chlorobenzene            | 6,200   | U      | 1,1,2,2-tetrachloroethane | 6,200   | U      |
| chloroethane             | 12,000  | U      | tetrachloroethene         | 6,200   | U      |
| 2-chloroethylvinyl ether | 12,000  | U      | toluene                   | 6,200   | U      |
| chloroform               | 6,200   | U      | 1,1,1-trichloroethane     | 6,200   | U      |
| chloromethane            | 12,000  | 'U     | 1,1,2-trichloroethane     | 6,200   | U      |
| dibromochloromethane     | 6,200   | U      | trichloroethene           | 6,200   | U      |
| 1,1-dichloroethane       | 6,200   | U      | trichlorofluoromethane    | 6,200   | U      |
| 1,2-dichloroethane       | 6,200   | Ü      | vinyl chloride            | 12,000  | Ü      |
|                          |         |        |                           |         |        |

Date of Analysis: 03/31/92

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

Client Project ID: LE Carpenter

Job Number: ITAD 51019

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in  $\mu g/kg$  (ppb)

Sample Matrix: Soil

Client Sample ID: 1045-052-1 Lab Sample ID: SS8495

| Tentative Identification (1) | Concentration (2) |
|------------------------------|-------------------|
| octane, 4-methyl-            | 11,000            |
| xylenes (total)              | 570,000 *         |
| nonane, 2,5-dimethyl-        | 26,000            |
| cyclohexanemethanol          | 18,000            |
| cyclohexane, propyl-         | 40,000            |
| benzene, ethylmethyl-        | 22,000 Y          |
| nonane, 2-methyl-            | 35,000            |
| 1,10-decanediol              | 6.500             |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- Quantitation based on continuing calibration.
- Y Indistinguishable isomer in tentatively identified compounds.

Job Number: ITAD 51019

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g/kg$  (ppb) Sample Matrix: Soil

Client Sample ID: 1045-052-1RE

Lab Sample ID: SS8495

| Compound                 | Concentra | tion         | Compound                  | Concent | ration |
|--------------------------|-----------|--------------|---------------------------|---------|--------|
| acrolein                 | 12,000    | U            | 1,1-dichloroethene        | 6,200   | U      |
| acrylonitrile            | 12,000    | U ,          | trans-1,2-dichloroethene  | 6,200   | Ú      |
| benzene                  | 6,200     | Ū            | 1,2-dichloropropane       | 6,200   | U      |
| bromodichloromethane     | 6,200     | U .          | cis-1,3-dichloropropene   | 6,200   | Ü      |
| bromoform                | 6,200     | Ü            | trans-1,3-dichloropropene | 6,200   | U      |
| bromomethane             | 12,000    | U .          | ethyl benzene             | 140,000 |        |
| carbon tetrachloride     | 6,200     | U            | methylene chloride        | 6,200   | U      |
| chlorobenzene            | 6,200     | U .          | 1,1,2,2-tetrachloroethane | 6,200   | U      |
| chloroethane             | 12,000    | Ü .          | tetrachloroethene         | 6,200   | U      |
| 2-chloroethylvinyl ether | 12,000    | U            | toluene                   | 6,200   | U      |
| chloroform               | 6,200     | Ü            | 1,1,1-trichloroethane     | 6,200   | U .    |
| chloromethane            | 12,000    | <b>U</b> : ' | 1,1,2-trichloroethane     | 6,200   | U      |
| dibromochloromethane     | 6,200     | U            | trichloroethene           | 6,200   | U      |
| 1,1-dichloroethane       | 6,200     | U            | trichlorofluoromethane    | 6,200   | U      |
| 1,2-dichloroethane       | 6,200     | <b>U</b> :.  | vinyl chloride            | 12,000  | U      |

Date of Analysis: 03/31/92

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Job Number: ITAD 51019

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/kg (ppb)

Sample Matrix: Soil

Client Sample ID: 1045-052-1RE Lab Sample ID: SS8495

| Tentative Identification (1) | Concentration (2) |
|------------------------------|-------------------|
| octane, 4-methyl-            | 12,000            |
| xylenes (total)              | 580,000 *         |
| nonane, 2,5-dimethyl-        | 30,000            |
| undecanal                    | 20,000            |
| cyclohexane, propyl-         | 44,000            |
| benzene, ethylmethyl-        | 24,000 Y          |
| nonane, 2-methyl-            | 42,000            |
| benzene, 1,3,5-trimethyl-    | 12,000            |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- Quantitation based on continuing calibration.
- Y Indistinguishable isomer in tentatively identified compounds.

Job Number: ITAD 51019

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in μg/kg (ppb) Sample Matrix: Soil

Client Sample ID: 1045-052-2 Lab Sample ID: SS8496

| Compound                 | Concentration | Compound                  | Concentration |
|--------------------------|---------------|---------------------------|---------------|
| acrolein                 | 10 U          | 1,1-dichloroethene        | 5 U           |
| acrylonitrile            | 10 Ü          | trans-1,2-dichloroethene  | 5 U           |
| benzene                  | 5 Ų           | 1,2-dichloropropane       | 5 U           |
| bromodichloromethane     | 5 U           | cis-1,3-dichloropropene   | 5 U           |
| bromoform                | 5 U           | trans-1,3-dichloropropene | 5 U           |
| bromomethane             | 10 U          | ethyl benzene             | 5 U           |
| carbon tetrachloride     | 5 U           | methylene chloride        | 7 B           |
| chlorobenzene            | 5 U           | 1,1,2,2-tetrachloroethane | 5 U           |
| chloroethane             | 10 U          | tetrachloroethene         | 5 U           |
| 2-chloroethylvinyl ether | 10 U          | toluene                   | 5 U           |
| chloroform               | 5 U           | 1,1,1-trichloroethane     | 5 U           |
| chloromethane            | 10 U          | 1,1,2-trichloroethane     | 5 U           |
| dibromochloromethane     | , <b>5</b> U  | trichloroethene           | 5 U           |
| 1,1-dichloroethane       | 5 U           | trichlorofluoromethane    | 5 U           |
| 1,2-dichloroethane       | 5 U           | vinyl chloride            | 10 U          |

Date of Analysis: 04/02/92

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

Client Project ID: LE Carpenter

Job Number: ITAD 51019

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/kg (ppb)

Sample Matrix: Soil

Client Sample ID: 1045-052-2 Lab Sample ID: SS8496

| Tentative Identification (1)       | Concentration (2) |
|------------------------------------|-------------------|
| acetone                            | 3.7 *             |
| cyclohexane, trimethyl             | 9.8 Y             |
| xylene                             | 11 *              |
| cyclohexane, ethylmethyl-          | 35 Y              |
| cyclohexane, 2-propenyl-           | 15                |
| unknown (substituted cyclohexane)  | 31                |
| cyclohexane, (1,2-dimethylbutyl)-  | 53                |
| cyclohexane, 1,-ethyl-2,3-dimethyl | 40                |
| unknown (substituted cyclohexane)  | 35                |
| cyclohexane, 1,-methyl-2-propyl    | 28                |
| unknown                            | 21                |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- \* Quantitation based on continuing calibration.
- Y Indistinguishable isomer in tentatively identified compounds.

Job Number: ITAD 51019

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g/kg$  (ppb) Sample Matrix: Soil

Client Sample ID: 1045-052-3 Lab Sample ID: SS8497

| Compound                 | Concentration | Compound                  | Concent | ration |
|--------------------------|---------------|---------------------------|---------|--------|
| acrolein                 | 10 U          | 1,1-dichloroethene        | 5       | U      |
| acrylonitrile            | 10 U          | trans-1,2-dichloroethene  | 5       | U      |
| benzene                  | 5 U           | 1,2-dichloropropane       | 5       | U      |
| bromodichloromethane     | 5 U           | cis-1,3-dichloropropene   | 5       | U      |
| bromoform                | 5 U           | trans-1,3-dichloropropene | 5       | U      |
| bromomethane             | 10 U          | ethyl benzene             | 5       | U      |
| carbon tetrachloride     | 5 U           | methylene chloride        | 4       | JB     |
| chlorobenzene            | 5 U           | 1,1,2,2-tetrachloroethane | 5       | U      |
| chloroethane             | 10 U          | tetrachloroethene         | 5       | U      |
| 2-chloroethylvinyl ether | 10 U          | toluene                   | 5       | Ŭ      |
| chloroform               | 5 U           | 1,1,1-trichloroethane     | 5       | U .    |
| chloromethane            | 10 U          | 1,1,2-trichloroethane     | 5       | U      |
| dibromochloromethane     | 5 U           | trichloroethene           | 5       | U      |
| 1,1-dichloroethane       | 5 U           | trichlorofluoromethane    | 5       | U      |
| 1,2-dichloroethane       | 5 Ü           | vinyl chloride            | 10      | U      |

Date of Analysis: 04/02/92

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

Job Number: ITAD 51019

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in  $\mu g/kg$  (ppb)

Sample Matrix: Soil

Client Sample ID: 1045-052-3 Lab Sample ID: SS8497

Tentative Identification (1)

Concentration (2)

xylenes

2.2 \*

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- \* Quantitation based on continuing calibration.

Job Number: ITAD 51019

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g/kg$  (ppb) Sample Matrix: Soil

Client Sample ID: Method Blank 1

Lab Sample ID: VB0331

| Compound                 | Concentration   | Compound                  | Concentration |  |  |
|--------------------------|-----------------|---------------------------|---------------|--|--|
| acrolein                 | 1,300 U         | 1,1-dichloroethene        | 620 U         |  |  |
| acrylonitrile            | 1,300 U         | trans-1,2-dichloroethene  | 620 U         |  |  |
| benzene                  | 620 U           | 1,2-dichloropropane       | 620 U         |  |  |
| bromodichloromethane     | 620 U           | cis-1,3-dichloropropene   | 620 U         |  |  |
| bromoform                | 620 U           | trans-1,3-dichloropropene | <b>620</b> U  |  |  |
| bromomethane             | 1,300 U         | ethyl benzene             | 620 U         |  |  |
| carbon tetrachloride     | 620 U           | methylene chloride        | 120 J         |  |  |
| chlorobenzene            | 620 U           | 1,1,2,2-tetrachloroethane | <b>620</b> U  |  |  |
| chloroethane             | 1,300 U         | tetrachloroethene         | 620 U         |  |  |
| 2-chloroethylvinyl ether | 1, <b>300</b> U | toluene                   | <b>620</b> U  |  |  |
| chloroform               | 620 U           | 1,1,1-trichloroethane     | 620 U         |  |  |
| chloromethane            | 1,300 U         | 1,1,2-trichloroethane     | 620 U         |  |  |
| dibromochloromethane     | 620 U           | trichloroethene           | 620 U         |  |  |
| 1,1-dichloroethane       | 6 <b>20</b> U   | trichlorofluoromethane    | 620 U         |  |  |
| 1,2-dichloroethane       | 620 U           | vinyl chloride            | 1,300 U       |  |  |

Date of Analysis: 03/31/92

This method blank applies to the following samples: 1045-052-1 and 1045-052-1 RE.

RE = Reanalyzed

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

Job Number: ITAD 51019

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in  $\mu g/kg$  (ppb)

Sample Matrix: Soil

Client Sample ID: Method Blank 1

Lab Sample ID: VB0331

Tentative Identification (1)

Concentration (2)

acetone

200

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITAD 51019

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g/kg$  (ppb) Sample Matrix: Soil

Client Sample ID: Method Blank 2

Lab Sample ID: EB04022

| Compound                 | Concentration | Compound                  | Concent | ration |
|--------------------------|---------------|---------------------------|---------|--------|
| acrolein                 | 10 U          | 1,1-dichloroethene        | 5       | U      |
| acrylonitrile            | 10 U          | trans-1,2-dichloroethene  | 5       | U      |
| benzene                  | 5 U           | 1,2-dichloropropane       | 5       | U      |
| bromodichloromethane     | 5 U           | cis-1,3-dichloropropene   | 5       | U      |
| bromoform                | 5 U           | trans-1,3-dichloropropene | 5       | U      |
| bromomethane             | 10 U          | ethyl benzene             | 5       | U      |
| carbon tetrachloride     | 5 U           | methylene chloride        | 1       | J      |
| chlorobenzene            | 5 U           | 1,1,2,2-tetrachloroethane | 5       | U      |
| chloroethane             | 10 U          | tetrachloroethene         | 5       | U      |
| 2-chloroethylvinyl ether | 10 Ù          | toluene                   | 5       | U      |
| chloroform               | 5 U           | 1,1,1-trichloroethane     | 5       | U      |
| chloromethane            | 10 U          | 1,1,2-trichloroethane     | 5       | U      |
| dibromochloromethane     | 5 U           | trichloroethene           | 5       | U      |
| 1,1-dichloroethane       | <b>5</b> U    | trichlorofluoromethane    | 5       | U      |
| 1,2-dichloroethane       | 5 U           | vinyl chloride            | ÍÔ      | U.     |

Date of Analysis: 04/02/92

This method blank applies to the following samples: 1045-052-2 and 1045-052-3.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

Job Number: ITAD 51019

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/kg (ppb)

Sample Matrix: Soil

Client Sample ID: Method Blank 2 Lab Sample ID: EB04022

Client Project ID: LE Carpenter

Tentative Identification (1)

Concentration (2)

acetonitrile

15

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

Job Number: ITAD 51019

Client Project ID: LE Carpenter

#### SOIL SURROGATE PERCENT RECOVERY SUMMARY

#### **VOLATILE**

| Client Sample ID | Toluene-D8<br>(81-117%)* | BFB<br><u>(74-121%)*</u> | 1,2 Dichloroethane-D4<br>(70-121%)* |
|------------------|--------------------------|--------------------------|-------------------------------------|
| 1045-052-1       | 48 **                    | 78                       | 48 **                               |
| 1045-052-1RE     | 48 **                    | 75                       | 49 **                               |
| Method Blank 1   | <b>9</b> 6               | 96                       | 94                                  |
| 1045-052-2       | 117                      | 97                       | 100                                 |
| 1045-052-3       | 115                      | 92                       | 99                                  |
| Method Blank 2   | 98                       | 98                       | 99                                  |

RE = Reanalysis

<sup>\*</sup>Values in parenthesis represent contract required QC limits.

<sup>\*\*</sup>Values are outside of contract required QC limits.

Job Number: ITAD 51019

### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in  $\mu g/kg$  (ppb)

Sample Matrix: Soil

Client Sample ID: 1045-052-1 Lab Sample ID: SS8498

| Compound                   | Concentration | Compound                    | Concentration | <u>on</u> |
|----------------------------|---------------|-----------------------------|---------------|-----------|
| acenaphthene               | 1,800 U       | bis(2-chloroisopropyl)ether | 1,800 U       | J         |
| acenaphthylene             | 1,800 U       | bis(2-ethylhexyl)phthalate  | 1,900,000 E   | BD        |
| anthracene                 | 1,800 U       | 4-bromophenyl phenyl ether  | 1,800 U       | J         |
| benzidine                  | 9,000 U       | 2-chloronaphthalene         | 1,800 U       | J         |
| benzo(a)anthracene         | 1,800 U       | 4-chlorophenyl phenyl ether | 1,800 U       | J         |
| benzo(b)fluoranthene       | 1,800 U       | chrysene                    | 1,800 U       | J         |
| benzo(k)fluoranthene       | 1,800 U       | dibenz(a,h)anthracene       | 1,800 U       | J         |
| benzo(a)pyrene             | 1,800 U       | di-n-butylphthalate         | 11,000        |           |
| benzo(g,h,i)perylene       | 1,800 U       | 1,2-dichlorobenzene         | 1,800 U       | j         |
| butylbenzylphthalate       | 1,600 J       | 1,3-dichlorobenzene         | 1,800 U       | J         |
| bis(2-chloroethoxy)methane | 1,800 U       | 1,4-dichlorobenzene         | 1,800 U       | j.        |
| bis(2-chloroethyl)ether    | 1,800 U       |                             |               |           |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 03/30/92 Date of Analysis: 04/13/92

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

E - Compound exceeded calibration range of instrument.

D - Compound analyzed at a secondary dilution factor.

Job Number: ITAD 51019

### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/kg (ppb)

Sample Matrix: Soil

Client Sample ID: 1045-052-1 Lab Sample ID: SS8498

| Compound                  | Concentra | <u>ition</u>         | Compound                   | Concentra | tion |
|---------------------------|-----------|----------------------|----------------------------|-----------|------|
| 3,3'dichlorobenzidine     | 3,600     | U                    | hexachloroethane           | 1,800     | U    |
| diethylphthalate          | 1,600     | J                    | indeno(1,2,3-cd)pyrene     | 1,800     | U    |
| dimethyl phthalate        | 380       | Ĵ                    | isophorone                 | 1,800     | U    |
| 2,4-dinitrotoluene        | 1,800     | U                    | naphthalene                | 1,800     | U    |
| 2,6-dinitrotoluene        | 1,800     | U                    | nitrobenzene               | 1,800     | U    |
| di-n-octyl phthalate      | 4,700     |                      | n-nitroso-di-n-propylamine | 1,800     | U    |
| 1,2-diphenylhydrazine(1)  | 1,800     | U                    | n-nitrosodimethylamine     | 1,800     | U    |
| fluoranthene              | 1,800     | U                    | n-nitrosodiphenylamine(2)  | 1,800     | U    |
| fluorene                  | 1,800     | Ų                    | phenanthrene               | 1,800     | U    |
| hexachlorobenzene         | 1,800     | $\mathbf{U}_{\perp}$ | pyrene                     | 1,800     | U    |
| hexachlorobutadiene       | 1,800     | U                    | 1,2,4-trichlorobenzene     | 1,800     | U    |
| hexachlorocyclopentadiene | 1,800     | U                    |                            |           |      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 03/30/92 Date of Analysis: 04/13/92

J - Indicates an estimated value less than the detection limit.

<sup>(1) -</sup> Screened for as azobenzene

<sup>(2) -</sup> Detected as diphenylamine

Job Number: ITAD 51019

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in  $\mu g/kg$  (ppb) Sample Matrix: Soil

Client Sample ID: 1045-052-1 Lab Sample ID: SS8498

| Tentative Identification (1)                           | Concentration (2) |
|--|-------------------|
| 2-pentanone, 4-hydroxy-4-methyl-                       | 14,000 ÅB         |
| unknown ( <sup>C</sup> 10 alkane)                      | 8,600             |
| unknown (alkane)                                       | 7,300             |
| unknown (alkane)                                       | 30,000            |
| unknown (alkane)                                       | 8,700             |
| unknown (alkane)                                       | 6,100             |
| unknown (alkane)                                       | 6,900             |
| unknown (alkane)                                       | 5,800             |
| unknown (alkane)                                       | 11,000            |
| propanoic acid, 2-methyl-, 1-(1,1-dimethylethyl)-2-met | 16,000            |
| hexadecanoic acid                                      | 6,700             |
| unknown  | 110,000           |
| octadecanoic acid, 2-methylpropyl ester                | 9,600             |
| phosphoric acid, 2-ethylhexyl diphenyl ester           | 55,000            |
| unknown (alkane)                                       | 190,000           |
| unknown (phthalate)                                    | 12,000            |
| unknown (phthalate)                                    | 31,000            |
| unknown (phthalate)                                    | 16,000            |
| unknown  | 5,600             |
| unknown (phthalate)                                    | 10,000            |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- A Suspected aldol condensation product.
- B Analyte was found in the blank as well as the sample.

Job Number: ITAD 51019

#### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in  $\mu g/kg$  (ppb)

Sample Matrix: Soil

Client Sample ID: 1045-052-2 Lab Sample ID: SS8499

| Compound                   | Concent | tration | Compound                    | Concenti | ation |
|----------------------------|---------|---------|-----------------------------|----------|-------|
| acenaphthene               | 2,300   | U       | bis(2-chloroisopropyl)ether | 2,300    | U     |
| acenaphthylene             | 2,300   | U       | bis(2-ethylhexyl)phthalate  | 650,000  | BD    |
| anthracene                 | 2,300   | U       | 4-bromophenyl phenyl ether  | 2,300    | U     |
| benzidine                  | 11,000  | U       | 2-chloronaphthalene         | 2,300    | U     |
| benzo(a)anthracene         | 2,300   | U       | 4-chlorophenyl phenyl ether | 2,300    | U     |
| benzo(b)fluoranthene       | 2,300   | U       | chrysene                    | 2,300    | U     |
| benzo(k)fluoranthene       | 2,300   | U       | dibenz(a,h)anthracene       | 2,300    | U     |
| benzo(a)pyrene             | 2,300   | U .     | di-n-butylphthalate         | 250      | J     |
| benzo(g,h,i)perylene       | 2,300   | U       | 1,2-dichlorobenzene         | 2,300    | U     |
| butylbenzylphthalate       | 380     | J       | 1,3-dichlorobenzene         | 2,300    | U     |
| bis(2-chloroethoxy)methane | 2,300   | U       | 1,4-dichlorobenzene         | 2,300    | Ü     |
| bis(2-chloroethyl)ether    | 2,300   | U       |                             |          |       |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction:

03/30/92

Date of Analysis:

04/13/92

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

D - Compound analyzed at a secondary dilution factor.

Job Number: ITAD 51019

## BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/kg (ppb)

Sample Matrix: Soil

Client Sample ID: 1045-052-2 Lab Sample ID: SS8499

| Compound                  | Concentra | <u>ttion</u> | Compound                   | Concentration |   |
|---------------------------|-----------|--------------|----------------------------|---------------|---|
| 3,3'dichlorobenzidine     | 4,500     | U            | hexachloroethane           | 2,300         | U |
| diethylphthalate          | 2,300     | U            | indeno(1,2,3-cd)pyrene     | 2,300         | Ū |
| dimethyl phthalate        | 2,300     | U            | isophorone                 | 2,300         | U |
| 2,4-dinitrotoluene        | 2,300     | U            | naphthalene                | 2,300         | U |
| 2,6-dinitrotoluene        | 2,300     | Ü            | nitrobenzene               | 2,300         | U |
| di-n-octyl phthalate      | 1,400     | <b>J</b> ,   | n-nitroso-di-n-propylamine | 2,300         | U |
| 1,2-diphenylhydrazine(1)  | 2,300     | U            | n-nitrosodimethylamine     | 2,300         | U |
| fluoranthene              | 2,300     | U            | n-nitrosodiphenylamine(2)  | 2,300         | U |
| fluorene                  | 2,300     | U            | phenanthrene               | 2,300         | U |
| hexachlorobenzene         | 2,300     | U            | pyrene                     | 2,300         | U |
| hexachlorobutadiene       | 2,300     | U            | 1,2,4-trichlorobenzene     | 2,300         | U |
| hexachlorocyclopentadiene | 2,300     | U            |                            | 4             | 1 |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 03/30/92 Date of Analysis: 04/13/92

J - Indicates an estimated value less than the detection limit.

<sup>(1) -</sup> Screened for as azobenzene

<sup>(2) -</sup> Detected as diphenylamine

Job Number: ITAD 51019

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in  $\mu g/kg$  (ppb) Sample Matrix: Soil

Client Sample ID: 1045-052-2 Lab Sample ID: SS8499

|  | •         |          |
|--|-----------|----------|
| Tentative Identification (1)                           | Concentra | tion (2) |
| 2-pentanone, 4-hydroxy-4-methyl-                       | 14,000    | AB       |
| unknown (aldol?)                                       | 1,100     | A        |
| propanoic acid, 2-methyl-, 1-(1,1-dimethylethyl)-2-met | 22,000    |          |
| hexadecanoic acid                                      | 1,000     |          |
| unknown (amide)  | 1,200     |          |
| phosphoric acid, 2-ethylhexyl diphenyl ester           | 8,400     |          |
| unknown  | 4,200     |          |
| unknown (phthalate)                                    | 1,400     |          |
| unknown (phthalate)                                    | 4,000     |          |
| unknown (phthalate)                                    | 3,100     |          |
| unknown (phthalate)                                    | 2,000     |          |
| unknown (alkane)                                       | 1,500     |          |
| unknown (phthalate)                                    | 1,800     |          |
| unknown (alkane)                                       | 1,400     |          |
| unknown  | 1,200     |          |
| unknown  | 1,200     | _        |
| unknown (phthalate)                                    | 3,200     |          |
| unknown (phthalate)                                    | 1,300     |          |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- A Suspected aldol condensation product.
- B Analyte was found in the blank as well as the sample.

Job Number: ITAD 51019

#### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in  $\mu g/kg$  (ppb)

Sample Matrix: Soil

Client Sample ID: 1045-052-3 Lab Sample ID: SS8500

| Compound                   | Concen | tration | Compound                    | Concentr | ation |
|----------------------------|--------|---------|-----------------------------|----------|-------|
| acenaphthene               | 1,900  | Ü .     | bis(2-chloroisopropyl)ether | 1,900    | U     |
| acenaphthylene             | 1,900  | U .     | bis(2-ethylhexyl)phthalate  | 840,000  | BD    |
| anthracene                 | 340    | J       | 4-bromophenyl phenyl ether  | . 1,900  | Ù     |
| benzidine                  | 9,400  | Ü       | 2-chloronaphthalene         | 1,900    | U     |
| benzo(a)anthracene         | 850    | J       | 4-chlorophenyl phenyl ether | 1,900    | U     |
| benzo(b)fluoranthene       | 640    | J       | chrysene                    | 890      | J     |
| benzo(k)fluoranthene       | 770    | J       | dibenz(a,h)anthracene       | 250      | J.    |
| benzo(a)pyrene             | 720    | J '     | di-n-butylphthalate         | 31,000   | D     |
| benzo(g,h,i)perylene       | 1,900  | U       | 1,2-dichlorobenzene         | 1,900    | U     |
| butylbenzylphthalate       | 1,500  | J       | 1,3-dichlorobenzene         | 1,900    | U     |
| bis(2-chloroethoxy)methane | 1,900  | U       | 1,4-dichlorobenzene         | 1,900    | U     |
| bis(2-chloroethyl)ether    | 1,900  | U       |                             |          |       |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction:

03/30/92

Date of Analysis:

04/13/92

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

D - Compound analyzed at a secondary dilution factor.

Job Number: ITAD 51019

### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/kg (ppb)

Sample Matrix: Soil

Client Sample ID: 1045-052-3 Lab Sample ID: SS8500

| Compound                  | Concentration | Compound                   | Concentration   |  |  |
|---------------------------|---------------|----------------------------|-----------------|--|--|
| 3,3'dichlorobenzidine     | 3,800 U       | hexachloroethane           | 1,900 U         |  |  |
| diethylphthalate          | 3,700         | indeno(1,2,3-cd)pyrene     | 380 J           |  |  |
| dimethyl phthalate        | 760 J         | isophorone                 | 1, <b>900</b> U |  |  |
| 2,4-dinitrotoluene        | 1,900 U       | naphthalene                | 230 J           |  |  |
| 2,6-dinitrotoluene        | 1,900 U       | nitrobenzene               | 1,900 U         |  |  |
| di-n-octyl phthalate      | 3,000         | n-nitroso-di-n-propylamine | 1,900 U         |  |  |
| 1,2-diphenylhydrazine(1)  | 1,900 U       | n-nitrosodimethylamine     | 1,900 U         |  |  |
| fluoranthene              | 1,500 J       | n-nitrosodiphenylamine(2)  | 1, <b>900</b> U |  |  |
| fluorene                  | 1,900 U       | phenanthrene               | 1,000 J         |  |  |
| hexachlorobenzene         | 1,900 U       | pyrene                     | 1,200 J         |  |  |
| hexachlorobutadiene       | 1,900 U       | 1,2,4-trichlorobenzene     | 1,900 U         |  |  |
| hexachlorocyclopentadiene | 1,900 U       |                            |                 |  |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 03/30/92 Date of Analysis: 04/13/92

J - Indicates an estimated value less than the detection limit.

<sup>(1) -</sup> Screened for as azobenzene

<sup>(2) -</sup> Detected as diphenylamine

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

Client Project ID: LE Carpenter

Job Number: ITAD 51019

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in  $\mu g/kg$  (ppb) Sample Matrix: Soil

Client Sample ID: 1045-052-3 Lab Sample ID: SS8500

| Tentative Identification (1)                           | Concentra | tion (2) |
|--|-----------|----------|
| 2-pentanone, 4-hydroxy-4-methyl-                       | 14,000    | AB       |
| unknown (alkane)                                       | 3,800     |          |
| unknown  | 3,700     |          |
| propanoic acid, 2-methyl-, 1-(1,1-dimethylethyl)-2-met | 5,700     |          |
| hexadecanoic acid                                      | 5,700     |          |
| unknown  | 100,000   |          |
| octadecanoic acid, 2-methylpropyl ester                | 9,000     | ••       |
| phosphoric acid, 2-ethylhexyl diphenyl ester           | 56,000    |          |
| unknown (phthalate)                                    | 4,700     |          |
| unknown (alkane)                                       | 220,000   |          |
| unknown (phthalate)                                    | 19,000    |          |
| unknown (phthalate)                                    | 39,000    |          |
| unknown (phthalate)                                    | 4,600     |          |
| unknown (phthalate)                                    | 22,000    |          |
| unknown (phthalate)                                    | 33,000    |          |
| unknown (phthalate)                                    | 3,900     |          |
| unknown  | 3,600     |          |
| unknown (phthalate)                                    | 6,300     |          |
| unknown (alkane)                                       | 10,000    |          |
| unknown (phthalate)                                    | 4,500     |          |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- A Suspected aldol condensation product.
- B Analyte was found in the blank as well as the sample.

Job Number: ITAD 51019

### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/kg (ppb)

Sample Matrix: Soil

Client Sample ID: Method Blank 1

Lab Sample ID: BL0163

| Compound                   | Concen | tration  | Compound                    | Concenti | ration |
|----------------------------|--------|----------|-----------------------------|----------|--------|
| acenaphthene               | 330    | U        | bis(2-chloroisopropyl)ether | 330      | U      |
| acenaphthylene             | 330    | U        | bis(2-ethylhexyl)phthalate  | 36       | J      |
| anthracene                 | 330    | U        | 4-bromophenyl phenyl ether  | 330      | Ŭ      |
| benzidine                  | 1,700  | U        | 2-chloronaphthalene         | 330      | U      |
| benzo(a)anthracene         | 330    | U        | 4-chlorophenyl phenyl ether | 330      | U      |
| benzo(b)fluoranthene       | 330    | U        | chrysene                    | 330      | U      |
| benzo(k)fluoranthene       | 330    | U        | dibenz(a,h)anthracene       | 330      | U      |
| benzo(a)pyrene             | 330    | U        | di-n-butylphthalate         | 330      | U      |
| benzo(g,h,i)perylene       | 330    | U        | 1,2-dichlorobenzene         | 330      | U      |
| butylbenzylphthalate       | 330    | <b>U</b> | 1,3-dichlorobenzene         | 330      | U      |
| bis(2-chloroethoxy)methane | 330    | U        | 1,4-dichlorobenzene         | 330      | U      |
| bis(2-chloroethyl)ether    | 330    | U        |                             |          |        |

Date of Extraction: 03/30/92 Date of Analysis: 04/10/92

This method blank applies to the following samples: 1045-052-1, 1045-052-1 DL, 1045-052-2, 1045-052-2 DL, 1045-052-3, 1045-052-3 DL and 1045-052-3 DL2.

DL = Dilution

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Job Number: ITAD 51019

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in  $\mu g/kg$  (ppb)

Sample Matrix: Soil

Client Sample ID: Method Blank 1

Lab Sample ID: BL0163

| Concentration | Compound  | Concentration  |
|---------------|---|--|
| 660 U         | hexachloroethane  | 330 U  |
| 330 Ù         | indeno(1,2,3-cd)pyrene  | 330 U  |
| 330 U         | isophorone  | 330 U  |
| 330 U         | naphthalene   | 330 Ù  |
| 330 U         | nitrobenzene  | <b>330</b> U   |
| 330 U         | n-nitroso-di-n-propylamine  | 330 U  |
| 330 U         | n-nitrosodimethylamine  | 330 U  |
| 330 U         | n-nitrosodiphenylamine(2)   | 330 U  |
| 330 U         | phenanthrene  | 330 U  |
| 330 U         | pyrene  | 330 U  |
| 330 U         | 1,2,4-trichlorobenzene  | 330 U  |
| 330 U         |   |  |
|               | 330 U | 660 U hexachloroethane  330 U indeno(1,2,3-cd)pyrene  330 U isophorone  330 U naphthalene  330 U nitrobenzene  330 U n-nitroso-di-n-propylamine  330 U n-nitrosodimethylamine  330 U n-nitrosodiphenylamine(2)  330 U phenanthrene  330 U pyrene  330 U 1,2,4-trichlorobenzene |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

<sup>(1) -</sup> Screened for as azobenzene

<sup>(2) -</sup> Detected as diphenylamine

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

Job Number: ITAD 51019

Client Project ID: LE Carpenter

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in μg/kg (ppb) Sample Matrix: Soil

Client Sample ID: Method Blank 1

Lab Sample ID: BL0163

| Tentative Identification (1)     | Concentra | tion (2) |
|----------------------------------|-----------|----------|
| 3-penten-2-one, 4-methyl-        | 160       | A        |
| unknown (aldol?)                 | 390       | A        |
| 2-pentanone, 4-hydroxy-4-methyl- | 6,900     | A        |
| unknown                          | 550       |          |
| 5-hexen-2-one, 5-methyl-         | 280       | A        |
| unknown (halogenated benzene)    | 230       |          |
| unknown (amide)                  | 3,200     |          |

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

A - Suspected aldol condensation product.

Job Number: ITAD 51019

#### SOIL SURROGATE PERCENT RECOVERY SUMMARY

#### **SEMIVOLATILE**

| Client Sample ID | Nitro-<br>Benzene-D5<br>(23-120%)* | 2-Fluoro-<br>Biphenyl<br>(30-115%)* | Terphenyl-<br>D14<br>(18-137%)* | Phenol-D5 (24-113%)* | 2-Fluoro-<br>Phenol<br>(25-121%)* | 2,4,6-<br>Tribromo-<br>Phenol<br>(19-122%)* |
|------------------|------------------------------------|-------------------------------------|---------------------------------|----------------------|-----------------------------------|---|
| 1045-052-1       | 86                                 | 84                                  | 82                              | 74                   | 82                                | 86  |
| 1045-052-1 DL    | D                                  | D                                   | D                               | D                    | D                                 | D   |
| 1045-052-2       | 80                                 | 76                                  | 71                              | 70                   | 77                                | 77  |
| 1045-052-2 DL    | D                                  | D                                   | D                               | D                    | D                                 | . <b>D</b>                                  |
| 1045-052-3       | 100                                | 86                                  | 82                              | 80                   | 92                                | 85  |
| 1045-052-3 DL    | 63                                 | 62                                  | <b>5</b> 6                      |                      | 61                                | 55  |
| 1045-052-3 DL2   | D                                  | <b>D</b>                            | D                               | ď                    | Ď                                 | D   |
| Method Blank 1   | 88                                 | 88                                  | 92                              | 78                   | 73                                | 72  |

<sup>\*</sup>Values in parenthesis represent contract required QC limits.

D - Surrogates diluted out

DL - Dilution

Job Number: ITPK 51491

# VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS Results in μg/kg (ppb) dry weight

Sample Matrix: Soil

Client Sample ID: 1045-067-01 Lab Sample ID: K3220

| Compound                 | Concent | ration | Compound                  | Concent | ration |
|--------------------------|---------|--------|---------------------------|---------|--------|
| acrolein                 | 1,400   | U      | 1,1-dichloroethene        | 700     | Ú      |
| acrylonitrile            | 1,400   | Ū      | cis-1,2-dichloroethene    | 700     | U      |
| benzene                  | 700     | U      | trans-1,2-dichloroethene  | 700     | U      |
| bromodichloromethane     | 700     | U      | 1,2-dichloropropane       | 700     | U      |
| bromoform                | 700     | U      | cis-1,3-dichloropropene   | 700     | Ü      |
| bromomethane             | 1,400   | U      | trans-1,3-dichloropropene | 700     | U      |
| carbon tetrachloride     | 700     | U      | ethyl benzene             | 29,000  | D      |
| chlorobenzene            | 700     | U      | methylene chloride        | 820     | В      |
| chloroethane             | 1,400   | U      | 1,1,2,2-tetrachloroethane | 700     | Ü      |
| 2-chloroethylvinyl ether | 1,400   | U      | tetrachloroethene         | 700     | U      |
| chloroform               | 700     | U      | toluene                   | 700     | U      |
| chloromethane            | 1,400   | U      | 1,1,1-trichloroethane     | 700     | U      |
| dibromochloromethane     | 700     | U      | 1,1,2-trichloroethane     | 700     | U      |
| 1,1-dichloroethane       | 700     | U      | trichloroethene           | 700     | Ü      |
| 1,2-dichloroethane       | 700     | U      | trichlorofluoromethane    | 700     | U      |
|                          |         |        | vinyl chloride            | 1,400   | U      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/09/92

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

D - Compound analyzed at a secondary dilution factor.

Job Number: ITPK 51491

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in  $\mu g/kg$  (ppb) dry weight

Sample Matrix: Soil

Client Sample ID: 1045-067-01 Lab Sample ID: K3220

| Tentative Identification (1) | Concentration (2) |
|------------------------------|-------------------|
| benzene, dimethyl-           | 110,000 Y         |
| unknown hydrocarbon (cyclic? | 30,000            |
| unknown cyclic alkane (C10)  | 18,000            |
| unknown alkane (C10)         | 15,000            |
| unknown alkane (C10)         | 36,000            |
| unknown alkane (C10)         | 18,000            |
| unknown alkane (cyclic?)     | 21,000            |
| unknown hydrocarbon          | 28,000            |
| unknown                      | 17,000            |
| unknown alkane (C11)         | 17,000            |

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Y - Indistinguishable isomer in tentatively identified compounds.

Job Number: ITPK 51491

### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS** Results in $\mu g/kg$ (ppb) dry weight

Sample Matrix: Soil

Client Sample ID: 1045-067-02 Lab Sample ID: K3221

| Compound                 | Concentration | Compound                  | Concent | ration |
|--------------------------|---------------|---------------------------|---------|--------|
| acrolein                 | 14 U          | 1,1-dichloroethene        | 7       | U      |
| acrylonitrile            | 14 U          | cis-1,2-dichloroethene    | 7       | U      |
| benzene                  | 7 U           | trans-1,2-dichloroethene  | 7       | U      |
| bromodichloromethane     | 7 U           | 1,2-dichloropropane       | 7       | U      |
| bromoform                | 7 U           | cis-1,3-dichloropropene   | 7       | U      |
| bromomethane             | 14 U          | trans-1,3-dichloropropene | 7       | U      |
| carbon tetrachloride     | , 7 U         | ethyl benzene             | 7       | U      |
| chlorobenzene            | 7 Û           | methylene chloride        | 430     | BD     |
| chloroethane             | 14 U          | 1,1,2,2-tetrachloroethane | 7       | .U     |
| 2-chloroethylvinyl ether | 14 U          | tetrachloroethene         | 7       | U      |
| chloroform               | 7 U           | toluene                   | 7       | U      |
| chloromethane            | 14 U          | 1,1,1-trichloroethane     | 7       | U      |
| dibromochloromethane     | 7 U           | 1,1,2-trichloroethane     | 7       | U      |
| 1,1-dichloroethane       | 7 U           | trichloroethene           | 7       | U      |
| 1,2-dichloroethane       | 7 U           | trichlorofluoromethane    | 7       | Ü      |
|                          |               | vinyl chloride            | 14      | U      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/10/92

<sup>-</sup> Indicates an estimated value less than the detection limit. J

В - Analyte was found in the blank as well as the sample.

Compound analyzed at a secondary dilution factor. D

Job Number: ITPK 51491

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/kg (ppb) dry weight

Sample Matrix: Soil

Client Sample ID: 1045-067-02 Lab Sample ID: K3221

| Tentative Identification (1) | Concentration (2) |
|------------------------------|-------------------|
| cyclohexane, butyl-          | 7.4               |
| cyclohexane, ethyl-methyl-   | 15 Y              |
| cyclohexane, ethyl-methyl-   | 22 Y              |
| unknown cyclic               | 16                |
| unknown cyclic               | 28                |
| cyclohexane, diethyl-        | 14                |
| cyclohexane, 1-ethyl-2,3-dim | <b>24</b>         |
| unknown                      | 23                |
| cyclohexane, 1-methyl-2-prop | 17                |

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Y - Indistinguishable isomer in tentatively identified compounds.

Job Number: ITPK 51491

# VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS Results in μg/kg (ppb) dry weight Sample Matrix: Soil

Client Sample ID: 1045-067-03 Lab Sample ID: K3222

| Compound                 | Concentration | Compound                  | Concent | ration   |
|--------------------------|---------------|---------------------------|---------|----------|
| acrolein                 | 14 ; U        | 1,1-dichloroethene        | 7       | U        |
| acrylonitrile            | 14 U          | cis-1,2-dichloroethene    | . 7     | U        |
| benzene                  | 7 · U         | trans-1,2-dichloroethene  | 7       | U        |
| bromodichloromethane     | 7 U           | 1,2-dichloropropane       | 7       | U        |
| bromoform                | 7 U           | cis-1,3-dichloropropene   | 7       | U        |
| bromomethane             | 14 U          | trans-1,3-dichloropropene | 7       | U        |
| carbon tetrachloride     | 7 U           | ethyl benzene             | 7       | U        |
| chlorobenzene            | , 7 U         | methylene chloride        | 79      | В        |
| chloroethane             | 14 U          | 1,1,2,2-tetrachloroethane | 7       | U        |
| 2-chloroethylvinyl ether | 14 U          | tetrachloroethene         | 7       | U        |
| chloroform               | 7 U           | toluene                   | 7       | <b>u</b> |
| chloromethane            | 14 U          | 1,1,1-trichloroethane     | 7       | U        |
| dibromochloromethane     | 7 U           | 1,1,2-trichloroethane     | . 7     | U        |
| 1,1-dichloroethane       | 7 U           | trichloroethene           | 7       | U        |
| 1,2-dichloroethane       | 7 U           | trichlorofluoromethane    | . 7     | U        |
|                          |               | vinyl chloride            | 14      | U        |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/10/92

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

Job Number: ITPK 51491

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in  $\mu g/kg$  (ppb) dry weight

Sample Matrix: Soil

Client Sample ID: 1045-067-03 Lab Sample ID: K3222

| Tentative Identification (1) | Concentration (2) |
|------------------------------|-------------------|
| acetone                      | 8.7               |
| cyclohexane, ethyl-methyl-   | 14 Y              |
| cyclohexane, ethyl-methyl-   | 16 Y              |
| unknown                      | 14                |
| cyclohexane, undecyl-        | 36                |
| cyclohexane, diethyl-        | 12                |
| cyclohexane, 1-ethyl-2,3-dim | 23                |
| unknown                      | 13                |
| cyclohexane, 1-methyl-2-prop | 14                |

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Y - Indistinguishable isomer in tentatively identified compounds.

Job Number: ITPK 51491

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in µg/kg (ppb) dry weight Sample Matrix: Soil

Client Sample ID: 1045-067-04 Lab Sample ID: K3223

| Compound                 | Concentrati | ion | Compound                  | Concent | ration |
|--------------------------|-------------|-----|---------------------------|---------|--------|
| acrolein                 | 13 U        | J   | 1,1-dichloroethene        | 6       | U      |
| acrylonitrile            | 13 U        | J   | cis-1,2-dichloroethene    | 6       | U      |
| benzene                  | 3 Ј         |     | trans-1,2-dichloroethene  | 6       | U      |
| bromodichloromethane     | 6 U         | J   | 1,2-dichloropropane       | 6       | U      |
| bromoform                | 6 U         | J   | cis-1,3-dichloropropene   | 6       | U      |
| bromomethane             | 13 U        | J   | trans-1,3-dichloropropene | 6       | U      |
| carbon tetrachloride     | 6 U         | J   | ethyl benzene             | 9       |        |
| chlorobenzene            | 6 U         | j . | methylene chloride        | 38      | В      |
| chloroethane             | 13 U        | J   | 1,1,2,2-tetrachloroethane | 6       | U      |
| 2-chloroethylvinyl ether | 13 U        | J   | tetrachloroethene         | 6       | U      |
| chloroform               | 6 U         | J   | toluene                   | 3       | J      |
| chloromethane            | 13 U        | J   | 1,1,1-trichloroethane     | 6       | U      |
| dibromochloromethane     | 6 U         | J   | 1,1,2-trichloroethane     | 6       | U      |
| 1,1-dichloroethane       | 6 U         | J   | trichloroethene           | , 6     | U      |
| 1,2-dichloroethane       | 6 U         | J   | trichlorofluoromethane    | . 6     | U      |
| •                        |             |     | vinyl chloride            | 13      | U      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/10/92

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

Client Project ID: LE Carpenter

Job Number: ITPK 51491

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in  $\mu g/kg$  (ppb) dry weight

Sample Matrix: Soil

Client Sample ID: 1045-067-04 Lab Sample ID: K3223

| Tentative Identification (1) | Concentration (2) |
|------------------------------|-------------------|
| acetone                      | . 11              |
| benzene, dimethyl            | 130 Y             |
| benzene, 1,2-dimethyl        | 12                |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- Y Indistinguishable isomer in tentatively identified compounds.

Job Number: ITPK 51491

### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g/kg$  (ppb) Sample Matrix: Soil

Client Sample ID: Method Blank 1 Lab Sample ID: WB06093

| Compound                 | Concentration   | Compound                  | Concent      | ration |
|--------------------------|-----------------|---------------------------|--------------|--------|
| acrolein                 | 1, <b>200</b> U | 1,1-dichloroethene        | 620          | U      |
| acrylonitrile            | 1,200 U         | cis-1,2-dichloroethene    | 620          | U      |
| benzene                  | 620 U           | trans-1,2-dichloroethene  | 620          | U      |
| bromodichloromethane     | 620 U           | 1,2-dichloropropane       | 620          | U      |
| bromoform                | 620 U           | cis-1,3-dichloropropene   | 620          | U      |
| bromomethane             | 1, <b>200</b> U | trans-1,3-dichloropropene | 620          | U      |
| carbon tetrachloride     | 620 U           | ethyl benzene             | 620          | U      |
| chlorobenzene            | 620 U           | methylene chloride        | 360          | J      |
| chloroethane             | 1,200 U         | 1,1,2,2-tetrachloroethane | 620          | U .    |
| 2-chloroethylvinyl ether | 1,200 U         | tetrachloroethene         | , <b>620</b> | U      |
| chloroform               | 620 Ü           | toluene                   | 620          | U      |
| chloromethane            | 1, <b>200</b> U | 1,1,1-trichloroethane     | 620          | Ü      |
| dibromochloromethane     | 620 U           | 1,1,2-trichloroethane     | 620          | U      |
| 1,1-dichloroethane       | 620 U           | trichloroethene           | 620          | U      |
| 1,2-dichloroethane       | 620 U           | trichlorofluoromethane    | 620          | U      |
|                          |                 | vinyl chloride            | 1,200        | U      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/09/92

This method blank applies to the following samples: 1045-067-01 and 1045-067-01 DL.

DL - Dilution.

J - Indicates an estimated value less than the detection limit.

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

Client Project ID: LE Carpenter

Job Number: ITPK 51491

### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/kg (ppb)

Sample Matrix: Soil

Client Sample ID: Method Blank 1 Lab Sample ID: WB06093

Tentative Identification (1)

Concentration (2)

None Detected

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITPK 51491

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g/kg$  (ppb) Sample Matrix: Soil

Client Sample ID: Method Blank 2 Lab Sample ID: VB0610

| Compound                 | Concentration | Compound                  | Concentration |
|--------------------------|---------------|---------------------------|---------------|
| acrolein                 | 10 U          | 1,1-dichloroethene        | 5 U           |
| acrylonitrile            | 10 U          | cis-1,2-dichloroethene    | 5 U           |
| benzene                  | <b>5</b> , U  | trans-1,2-dichloroethene  | . 5 U         |
| bromodichloromethane     | <b>5</b> U    | 1,2-dichloropropane       | 5 U           |
| bromoform                | <b>5</b> U    | cis-1,3-dichloropropene   | 5 U           |
| bromomethane             | 10 U          | trans-1,3-dichloropropene | 5 U           |
| carbon tetrachloride     | <b>5</b> U    | ethyl benzene             | 5 U           |
| chlorobenzene            | <b>5</b> U    | methylene chloride        | 2 J           |
| chloroethane             | 10 U          | 1,1,2,2-tetrachloroethane | 5 U           |
| 2-chloroethylvinyl ether | 10 , U        | tetrachloroethene         | 5 U           |
| chloroform               | 5 U           | toluene                   | 5 U           |
| chloromethane            | 10 U          | 1,1,1-trichloroethane     | 5 U           |
| dibromochloromethane     | <b>5 U</b>    | 1,1,2-trichloroethane     | . 5 U         |
| 1,1-dichloroethane       | 5 U           | trichloroethene           | 5 U           |
| 1,2-dichloroethane       | 5 U           | trichlorofluoromethane    | 5 U           |
|                          |               | vinyl chloride            | 10 U          |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/10/92

This method blank applies to the following samples: 1045-067-02, 1045-067-02 DL, 1045-067-03 and 1045-067-04.

DL - Dilution.

J - Indicates an estimated value less than the detection limit.

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

Client Project ID: LE Carpenter

Job Number: ITPK 51491

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/kg (ppb)

Sample Matrix: Soil

Client Sample ID: Method Blank 2

Lab Sample ID: VB0610

Tentative Identification (1)

Concentration (2)

None Detected

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITPK 51491

### SOIL SURROGATE PERCENT RECOVERY SUMMARY

|                  | •••                      | VOLATILE          |                                     |
|------------------|--------------------------|-------------------|-------------------------------------|
| Client Sample ID | Toluene-D8<br>(81-117%)* | BFB<br>(74-121%)* | 1,2 Dichloroethane-D4<br>(70-121%)* |
| 1045-067-01      | 97                       | 154 **            | 99                                  |
| 1045-067-01 DL   | 102                      | 110               | 91                                  |
| Method Blank 1   | 101                      | 98                | 94                                  |
| 1045-067-02      | 106                      | 109               | 93                                  |
| 1045-067-02 DL   | 105                      | 108               | 96                                  |
| 1045-067-03      | 105                      | 112               | 97                                  |
| 1045-067-04      | 105                      | 105               | 95                                  |
| Method Blank 2   | 104                      | 103               | 101                                 |

DL - Dilution

<sup>\*</sup>Values in parenthesis represent QC limits.

<sup>\*\*</sup>Values are outside of QC limits.

Job Number: ITPK 51491

#### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in  $\mu g/kg$  (ppb) dry weight

Sample Matrix: Soil

Client Sample ID: 1045-067-01 Lab Sample ID: K3228

| Compound                   | Concen | tration  | Compound                    | Concentr | ration |
|----------------------------|--------|----------|-----------------------------|----------|--------|
| acenaphthene               | 740    | U        | bis(2-chloroisopropyl)ether | 740      | U      |
| acenaphthylene             | 740    | Ŭ        | bis(2-ethylhexyl)phthalate  | 110,000  | D      |
| anthracene                 | 740    | U        | 4-bromophenyl phenyl ether  | 740      | Ü      |
| benzidine                  | 3,700  | U        | 2-chloronaphthalene         | 740      | Ų      |
| benzo(a)anthracene         | 740    | U        | 4-chlorophenyl phenyl ether | 740      | U      |
| benzo(b)fluoranthene       | 740    | U.       | chrysene                    | 740      | U      |
| benzo(k)fluoranthene       | 740    | U        | dibenz(a,h)anthracene       | 740      | U      |
| benzo(a)pyrene             | 740    | U        | di-n-butylphthalate         | 600      | J      |
| benzo(g,h,i)perylene       | 740    | U        | 1,2-dichlorobenzene         | 740      | U      |
| butylbenzylphthalate       | 1,800  | <b>B</b> | 1,3-dichlorobenzene         | 740      | U      |
| bis(2-chloroethoxy)methane | 740    | Ū        | 1,4-dichlorobenzene         | 740      | U      |
| bis(2-chloroethyl)ether    | 740    | Ŭ.       |                             |          |        |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

D - Compound analyzed at a secondary dilution factor.

Job Number: ITPK 51491

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

#### Results in µg/kg (ppb) dry weight

Sample Matrix: Soil

Client Sample ID: 1045-067-01 Lab Sample ID: K3228

| Compound                  | Concentra | tion | Compound                   | Concentra | ation |
|---------------------------|-----------|------|----------------------------|-----------|-------|
| 3,3'dichlorobenzidine     | 1,500     | U    | hexachloroethane           | 740       | Ų     |
| diethylphthalate          | 740       | U    | indeno(1,2,3-cd)pyrene     | 740       | U     |
| dimethyl phthalate        | 740       | U    | isophorone                 | 740       | U     |
| 2,4-dinitrotoluene        | 740       | Ü    | naphthalene                | 740       | U     |
| 2,6-dinitrotoluene        | 740       | Ŭ    | nitrobenzene               | 740       | U     |
| di-n-octyl phthalate      | 2,500     |      | n-nitroso-di-n-propylamine | 740       | U     |
| 1,2-diphenylhydrazine(1)  | 740       | U    | n-nitrosodimethylamine     | 740       | U     |
| fluoranthene              | 740       | U    | n-nitrosodiphenylamine(2)  | 740       | U     |
| fluorene                  | 740       | U    | phenanthrene               | 740       | U     |
| hexachlorobenzene         | 740       | Ŭ    | pyrene                     | 740       | U     |
| hexachlorobutadiene       | 740       | U    | 1,2,4-trichlorobenzene     | 740       | U     |
| hexachlorocyclopentadiene | 740       | U    |                            |           |       |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

(1) - Screened for as azobenzene

(2) - Detected as diphenylamine

Date of Extraction: 00
Date of Analysis: 00

06/02/92 06/08/92

Job Number: ITPK 51491

#### ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/kg (ppb) dry weight

Sample Matrix: Soil

Client Sample ID: 1045-067-01 Lab Sample ID: K3228

| Compound                   | Concentration | Compound              | Concentration |
|----------------------------|---------------|-----------------------|---------------|
| 4-chloro-3-methylphenol    | 740 U         | 2-nitrophenol         | 740 U         |
| 2-chlorophenol             | 740 U         | 4-nitrophenol         | 3,700 U       |
| 2,4-dichlorophenol         | 740 U         | pentachlorophenol     | 3,700 U       |
| 2,4-dimethylphenol         | 740 U         | phenol                | 740 U         |
| 2,4-dinitrophenol          | 3,700 U       | 2,4,6-trichlorophenol | 740 U         |
| 2-methyl-4,6-dinitrophenol | 3,700 U       |                       |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Job Number: ITPK 51491

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/kg (ppb) dry weight

Sample Matrix: Soil

Client Sample ID: 1045-067-01 Lab Sample ID: K3228

| Tentative Identification (1)     | Concentration (2) |
|----------------------------------|-------------------|
| unknown                          | 12,000            |
| ethanol, 2-(dodecyloxy)-         | 11,000            |
| cyclopropane, nonyl-             | 11,000            |
| unknown                          | 34,000            |
| unknown                          | 29,000            |
| 1-hexadecanol                    | 9,100             |
| ethanol, 2-(dodecyloxy)-         | 8,800             |
| unknown                          | 24,000            |
| 2-hexadecanol                    | 26,000            |
| unknown                          | 22,000            |
| 1-hexadecanol                    | 5,600             |
| hexanedioic acid, dioctyl ester  | 16,000 B          |
| unknown                          | 25,000            |
| 2-pentanone, 4-hydroxy-4-methyl- | 25,000 AB         |
| ethanol, 2-(hexadecyloxy)-       | 4,400             |
| unknown                          | 13,000            |
| 1,2-benzenedicarboxylic acid     | 13,000            |
| 1,2-benzenedicarboxylic acid     | 12,000            |
| unknown                          | 12,000            |
| octane, 2,3-dimethyl-            | 19,000            |
| propanoic acid, 2-methyl-, 1     | 3,100             |
| unknown                          | 17,000            |
| 1,2-benzenedicarboxylic acid     | 9,700             |
| octane, 3,6-dimethyl-            | 16,000            |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- A Suspected aldol condensation product.
- B Analyte was found in the blank as well as the sample.

Job Number: ITPK 51491

#### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/kg (ppb) dry weight

Sample Matrix: Soil

Client Sample ID: 1045-067-02 Lab Sample ID: K3229

| Compound                   | Concentration | Compound                    | Concentration |
|----------------------------|---------------|-----------------------------|---------------|
| acenaphthene               | 1,800 U       | bis(2-chloroisopropyl)ether | 1,800 U       |
| acenaphthylene             | 1,800 U       | bis(2-ethylhexyl)phthalate  | 20,000        |
| anthracene                 | 1,800 U       | 4-bromophenyl phenyl ether  | 1,800 U       |
| benzidine                  | 9,100 U       | 2-chloronaphthalene         | 1,800 U       |
| benzo(a)anthracene         | 1,800 U       | 4-chlorophenyl phenyl ether | 1,800 U       |
| benzo(b)fluoranthene       | 1,800 U       | chrysene                    | 1,800 U       |
| benzo(k)fluoranthene       | 1,800 U       | dibenz(a,h)anthracene       | 1,800 U       |
| benzo(a)pyrene             | 1,800 U       | di-n-butylphthalate         | 1,800 U       |
| benzo(g,h,i)perylene       | 1,800 U       | 1,2-dichlorobenzene         | 1,800 U       |
| butylbenzylphthalate       | 1,800 U       | 1,3-dichlorobenzene         | 1,800 U       |
| bis(2-chloroethoxy)methane | 1,800 U       | 1,4-dichlorobenzene         | 1,800 U       |
| bis(2-chloroethyl)ether    | 1,800 U       |                             |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Job Number: ITPK 51491

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

#### Results in $\mu g/kg$ (ppb) dry weight

Sample Matrix: Soil

Client Sample ID: 1045-067-02 Lab Sample ID: K3229

| Compound                  | Concentra | tion | Compound                   | Concentra | ation |
|---------------------------|-----------|------|----------------------------|-----------|-------|
| 3,3'dichlorobenzidine     | 3,600     | Ü    | hexachloroethane           | 1,800     | Ŭ     |
| diethylphthalate          | 1,800     | U    | indeno(1,2,3-cd)pyrene     | 1,800     | U     |
| dimethyl phthalate        | 1,800     | U    | isophorone                 | 1,800     | U     |
| 2,4-dinitrotoluene        | 1,800     | U    | naphthalene                | 1,800     | U     |
| 2,6-dinitrotoluene        | 1,800     | Ü    | nitrobenzene               | 1,800     | U     |
| di-n-octyl phthalate      | 310       | J    | n-nitroso-di-n-propylamine | 1,800     | U     |
| 1,2-diphenylhydrazine(1)  | 1,800     | U    | n-nitrosodimethylamine     | 1,800     | U     |
| fluoranthene              | 1,800     | U    | n-nitrosodiphenylamine(2)  | 1,800     | U     |
| fluorene                  | 1,800     | U    | phenanthrene               | 1,800     | U     |
| hexachlorobenzene         | 1,800     | U    | pyrene                     | 1,800     | U     |
| hexachlorobutadiene       | 1,800     | U    | 1,2,4-trichlorobenzene     | 1,800     | U     |
| hexachlorocyclopentadiene | 1,800     | Ü    |                            |           |       |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

(1) - Screened for as azobenzene

(2) - Detected as diphenylamine

Job Number: ITPK 51491

#### ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/kg (ppb) dry weight

Sample Matrix: Soil

Client Sample ID: 1045-067-02 Lab Sample ID: K3229

| Compound                   | Concentration | Compound              | Concentration |
|----------------------------|---------------|-----------------------|---------------|
| 4-chloro-3-methylphenol    | 1,800 U       | 2-nitrophenol         | 1,800 U       |
| 2-chlorophenol             | 1,800 U       | 4-nitrophenol         | 9,100 U       |
| 2,4-dichlorophenol         | 1,800 U       | pentachlorophenol     | 9,100 U       |
| 2,4-dimethylphenol         | 1,800 U       | phenol                | 1,800 U       |
| 2,4-dinitrophenol          | 9,100 U       | 2,4,6-trichlorophenol | 1,800 U       |
| 2-methyl-4,6-dinitrophenol | 9,100 U       |                       |               |

Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction:

06/02/92

Date of Analysis:

06/08/92

Indicates an estimated value less than the detection limit.

Job Number: ITPK 51491

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/kg (ppb) dry weight

Sample Matrix: Soil

Client Sample ID: 1045-067-02 Lab Sample ID: K3229

| Tentative Identification (1) Concentr |        | n (2) |
|---------------------------------------|--------|-------|
| hexanedioic acid, dioctyl ester       | 16,000 | B     |
| unknown                               | 16,000 |       |
| 2-pentanone, 4-hydroxy-4-methyl-      | 17,000 | AB    |
| 1,2-benzenedicarboxylic acid-         | 870    |       |
| 2H-pyran-2,3-diol, tetrahydr          | 1,500  |       |
| 2-pentanone,5-(acetyloxy)-            | 1,200  | ΑB    |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- A Suspected aldol condensation product.
- B Analyte was found in the blank as well as the sample.

Job Number: ITPK 51491

#### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/kg (ppb) dry weight

Sample Matrix: Soil

Client Sample ID: 1045-067-03 Lab Sample ID: K3230

| Compound                   | Concentrat | ncentration Compound |                             | Concentration |   |
|----------------------------|------------|----------------------|-----------------------------|---------------|---|
| acenaphthene               | 900 U      | J                    | bis(2-chloroisopropyl)ether | 900           | U |
| acenaphthylene             | 900 U      | J                    | bis(2-ethylhexyl)phthalate  | 11,000        |   |
| anthracene                 | 900 T      | J :                  | 4-bromophenyl phenyl ether  | 900           | U |
| benzidine                  | 4,500 U    | J                    | 2-chloronaphthalene         | 900           | U |
| benzo(a)anthracene         | 900 T      | J ·                  | 4-chlorophenyl phenyl ether | 900           | U |
| benzo(b)fluoranthene       | 900 U      | J                    | chrysene                    | 900           | U |
| benzo(k)fluoranthene       | 900 U      | J                    | dibenz(a,h)anthracene       | 900           | U |
| benzo(a)pyrene             | 120 J      | •                    | di-n-butylphthalate         | .900          | U |
| benzo(g,h,i)perylene       | 900 U      | J                    | 1,2-dichlorobenzene         | 900           | U |
| butylbenzylphthalate       | 900 U      | J                    | 1,3-dichlorobenzene         | 900           | U |
| bis(2-chloroethoxy)methane | 900 U      | J                    | 1,4-dichlorobenzene         | 900           | U |
| bis(2-chloroethyl)ether    | 900 T      | J.,                  |                             |               |   |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction:

06/02/92

Date of Analysis:

06/08/92

J - Indicates an estimated value less than the detection limit.

Job Number: ITPK 51491

#### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

#### Results in $\mu g/kg$ (ppb) dry weight

Sample Matrix: Soil

Client Sample ID: 1045-067-03 Lab Sample ID: K3230

| Compound                  | <b>Concentration</b> |     | Concentration Compound     |     | Compound | Concentration |  |  |
|---------------------------|----------------------|-----|----------------------------|-----|----------|---------------|--|--|
| 3,3'dichlorobenzidine     | 1,800                | U   | hexachloroethane           | 900 | Ŭ        |               |  |  |
| diethylphthalate          | 900                  | U   | indeno(1,2,3-cd)pyrene     | 900 | U        |               |  |  |
| dimethyl phthalate        | 900                  | U   | isophorone                 | 900 | U        |               |  |  |
| 2,4-dinitrotoluene        | 900                  | U   | naphthalene                | 900 | Ú        |               |  |  |
| 2,6-dinitrotoluene        | 900                  | U   | nitrobenzene               | 900 | Ų        |               |  |  |
| di-n-octyl phthalate      | 440                  | J ; | n-nitroso-di-n-propylamine | 900 | U        |               |  |  |
| 1,2-diphenylhydrazine(1)  | 900                  | U   | n-nitrosodimethylamine     | 900 | U        |               |  |  |
| fluoranthene              | 900                  | U   | n-nitrosodiphenylamine(2)  | 900 | U        |               |  |  |
| fluorene                  | 900                  | U   | phenanthrene               | 900 | Ū        |               |  |  |
| hexachlorobenzene         | 900                  | U   | pyrene                     | 900 | Ù        |               |  |  |
| hexachlorobutadiene       | 900                  | U   | 1,2,4-trichlorobenzene     | 900 | Ų        |               |  |  |
| hexachlorocyclopentadiene | 900                  | U   |                            |     |          |               |  |  |

- Compound was analyzed for but not detected. The number is the detection limit for the sample.

- Indicates an estimated value less than the detection limit.

(1) - Screened for as azobenzene

(2) - Detected as diphenylamine

Date of Extraction:

06/02/92

Date of Analysis:

06/08/92

Job Number: ITPK 51491

#### ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

### Results in $\mu g/kg$ (ppb) dry weight

Sample Matrix: Soil

Client Sample ID: 1045-067-03 Lab Sample ID: K3230

| Compound                   | <u>Concentration</u> <u>Compound</u> |   | Compound              | Concentration |   |
|----------------------------|--------------------------------------|---|-----------------------|---------------|---|
| 4-chloro-3-methylphenol    | 900                                  | U | 2-nitrophenol         | 900           | U |
| 2-chlorophenol             | 900                                  | U | 4-nitrophenol         | 4,500         | U |
| 2,4-dichlorophenol         | 900                                  | U | pentachlorophenol     | 4,500         | U |
| 2,4-dimethylphenol         | 900                                  | U | phenol                | 900           | U |
| 2,4-dinitrophenol          | 4,500                                | U | 2,4,6-trichlorophenol | 900           | Ú |
| 2-methyl-4,6-dinitrophenol | 4,500                                | U |                       |               |   |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

"lient Project ID: LE Carpenter

Job Number: ITPK 51491

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in  $\mu g/kg$  (ppb) dry weight

Sample Matrix: Soil

Client Sample ID: 1045-067-03 Lab Sample ID: K3230

| Tentative Identification (1)    | Concentration (2) |    |  |
|---------------------------------|-------------------|----|--|
| hexanedioic acid, dioctÿl ester | 8,100             | B  |  |
| unknown                         | 7,900             |    |  |
| 2-pentanone, 4-hydroxy-4-methyl | 12,000            | AB |  |
| 2-pentanone,5-(acetyloxy)-      | 1,100             | AB |  |
| phosphoric acid, 2-ethylhexy    | 370               |    |  |
| 2H-pyran-2,3-diol, tetrahydr    | 850               |    |  |
| 1,2-benzenedicarboxylic acid    | 410               |    |  |
| undecane, 5-methyl-             | 480               |    |  |
| 2-heptanol, acetate             | 470               | В  |  |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- A Suspected aldol condensation product.
- B Analyte was found in the blank as well as the sample.

Job Number: ITPK 51491

#### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

#### Results in $\mu g/kg$ (ppb) dry weight

Sample Matrix: Soil

Client Sample ID: 1045-067-04 Lab Sample ID: K3231

| Compound                   | Concentration    | Compound                    | Concentration   |
|----------------------------|------------------|-----------------------------|-----------------|
| acenaphthene               | 1,700 U          | bis(2-chloroisopropyl)ether | 1,700 U         |
| acenaphthylene             | 1,700 U          | bis(2-ethylhexyl)phthalate  | 310,000 D       |
| anthracene                 | 1,700 U          | 4-bromophenyl phenyl ether  | 1,700 U         |
| benzidine                  | 8,400 U          | 2-chloronaphthalene         | 1,700 U         |
| benzo(a)anthracene         | 1,700 U          | 4-chlorophenyl phenyl ether | 1,700 U         |
| benzo(b)fluoranthene       | 1,700 U          | chrysene                    | 1,700 U         |
| benzo(k)fluoranthene       | 1,700 U          | dibenz(a,h)anthracene       | 1,700 U         |
| benzo(a)pyrene             | 1, <b>70</b> 0 U | di-n-butylphthalate         | 1, <b>700</b> U |
| benzo(g,h,i)perylene       | 1,700 U          | 1,2-dichlorobenzene         | 1,700 U         |
| butylbenzylphthalate       | 600 J            | 1,3-dichlorobenzene         | 1,700 U         |
| bis(2-chloroethoxy)methane | 1,700 U          | 1,4-dichlorobenzene         | 1,700 U         |
| bis(2-chloroethyl)ether    | 1, <b>700</b> U  |                             |                 |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

D - Compound analyzed at a secondary dilution factor.

Job Number: ITPK 51491

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

#### Results in $\mu g/kg$ (ppb) dry weight

Sample Matrix: Soil

Client Sample ID: 1045-067-04 Lab Sample ID: K3231

| Compound                  | Concentration |    | Compound                   | Concentration |   |  |
|---------------------------|---------------|----|----------------------------|---------------|---|--|
| 3,3'dichlorobenzidine     | 3,400         | U  | hexachloroethane           | 1,700         | Ų |  |
| diethylphthalate          | 1,700         | U  | indeno(1,2,3-cd)pyrene     | 1,700         | U |  |
| dimethyl phthalate        | 1,700         | U  | isophorone                 | 1,700         | U |  |
| 2,4-dinitrotoluene        | 1,700         | U. | naphthalene                | 1,700         | U |  |
| 2,6-dinitrotoluene        | 1,700         | U  | nitrobenzene               | 1,700         | U |  |
| di-n-octyl phthalate      | 1,100         | Ĵ  | n-nitroso-di-n-propylamine | 1,700         | U |  |
| 1,2-diphenylhydrazine(1)  | 1,700         | U  | n-nitrosodimethylamine     | 1,700         | U |  |
| fluoranthene              | 1,700         | U  | n-nitrosodiphenylamine(2)  | 1,700         | U |  |
| fluorene                  | 1,700         | U  | phenanthrene               | 1,700         | U |  |
| hexachlorobenzene         | 1,700         | U  | pyrene                     | 1,700         | U |  |
| hexachlorobutadiene       | 1,700         | U  | 1,2,4-trichlorobenzene     | 1,700         | U |  |
| hexachlorocyclopentadiene | 1,700         | U  |                            | ·             |   |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

(1) - Screened for as azobenzene

(2) - Detected as diphenylamine

Date of Extraction:
Date of Analysis:

06/02/92 06/08/92

Job Number: ITPK 51491

#### ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in  $\mu g/kg$  (ppb) dry weight

Sample Matrix: Soil

Client Sample ID: 1045-067-04 Lab Sample ID: K3231

| Compound                   | Concentration   | Compound              | Concentration |  |
|----------------------------|-----------------|-----------------------|---------------|--|
| 4-chloro-3-methylphenol    | 1, <b>700</b> U | 2-nitrophenol         | 1,700 U       |  |
| 2-chlorophenol             | 1,700 U         | 4-nitrophenol         | 8,400 U       |  |
| 2,4-dichlorophenol         | 1, <b>700</b> U | pentachlorophenol     | 8,400 U       |  |
| 2,4-dimethylphenol         | 1,700 U         | phenol                | 1,700 U       |  |
| 2,4-dinitrophenol          | 8,400 U         | 2,4,6-trichlorophenol | 1,700 U       |  |
| 2-methyl-4,6-dinitrophenol | 8,400 Ü         |                       |               |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Job Number: ITPK 51491

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/kg (ppb) dry weight

Sample Matrix: Soil

Client Sample ID: 1045-067-04 Lab Sample ID: K3231

| Tentative Identification (1)      | Concentration (2) |
|-----------------------------------|-------------------|
| hexadecanoic                      | 14,000            |
| cyclooctane, -dimethyl-           | 1 <b>4,000</b> Y  |
| propanoic acid, 2-methyl-, 1      | 13,000            |
| hexanedioic acid, dioctyl ester . | 14,000 B          |
| 2-pentanone, 4-hydroxy-4-methyl-  | 15,000 AB         |
| tetradecanoic acid                | 6,500             |
| 1,2-benzenedicarboxylic acid      | 8,600             |
| phosphoric acid, 2-ethylhexy      | 7,400             |
| 1,2-benzenedicarboxylic acid      | 5,400             |
| unknown                           | 4,000             |
| dodecanamide, N,N-bis(2-hydr      | 2,100             |
| 1,2-benzenedicarboxylic acid      | 3,100             |
| 14-pentadecenoic acid             | 1,500             |
| unknown                           | 2,500             |
| 1,2-benzenedicarboxylic acid      | 1,700             |
| unknown                           | 2,000             |
| unknown                           | 1,600             |
| 1,2-benzenedicarboxylic acid      | 1,500             |
| propanoic acid, 2-methyl-, 2      | 810               |
| 2H-pyran-2,3-diol, tetrahydr      | 1,600             |
| benzene, ethyl methyl             | 1,600 Y           |
| benzene, methyl(-methylethyl)     | 1,300 Y           |
| benzene, trimethyl-               | 1,200 Y           |
| heptane, 2,4-dimethyl-            | 1,100             |
|                                   |                   |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- A Suspected aldol condensation product.
- B Analyte was found in the blank as well as the sample.
- Y Indistinguishable isomer in tentatively identified compounds.

Job Number: ITPK 51491

#### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in  $\mu g/kg$  (ppb)

Sample Matrix: Soil

Client Sample ID: Method Blank

Lab Sample ID: H0675

| Compound                   | Concentr | ation | Compound                    | Concentr | ation |
|----------------------------|----------|-------|-----------------------------|----------|-------|
| acenaphthene               | 330      | U     | bis(2-chloroisopropyl)ether | 330      | Ü     |
| acenaphthylene             | 330      | U     | bis(2-ethylhexyl)phthalate  | 330      | U     |
| anthracene                 | 330      | U     | 4-bromophenyl phenyl ether  | 330      | U     |
| benzidine                  | 1,700    | U     | 2-chloronaphthalene         | 330      | U     |
| benzo(a)anthracene         | 330      | U     | 4-chlorophenyl phenyl ether | 330      | U     |
| benzo(b)fluoranthene       | 330      | U     | chrysene                    | 330      | U     |
| benzo(k)fluoranthene       | 330      | U     | dibenz(a,h)anthracene       | 330      | U     |
| benzo(a)pyrene             | 330      | U     | di-n-butylphthalate         | 330      | U     |
| benzo(g,h,i)perylene       | 330      | U     | 1,2-dichlorobenzene         | 330      | U     |
| butylbenzylphthalate       | 48       | J     | 1,3-dichlorobenzene         | 330      | U     |
| bis(2-chloroethoxy)methane | 330      | U     | 1,4-dichlorobenzene         | 330      | U     |
| bis(2-chloroethyl)ether    | 330      | U     |                             |          |       |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction:

06/02/92

Date of Analysis:

06/08/92

J - Indicates an estimated value less than the detection limit.

Job Number: ITPK 51491

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/kg (ppb)

Sample Matrix: Soil

Client Sample ID: Method Blank Lab Sample ID: H0675

| Compound                  | Concentration |          | Compound                   | Concentration |   |  |
|---------------------------|---------------|----------|----------------------------|---------------|---|--|
| 3,3'dichlorobenzidine     | 660           | <b>U</b> | hexachloroethane           | 330           | U |  |
| diethylphthalate          | 330           | U        | indeno(1,2,3-cd)pyrene     | 330           | Ü |  |
| dimethyl phthalate        | 330           | Ü        | isophorone                 | 330           | U |  |
| 2,4-dinitrotoluene        | 330           | U        | naphthalene                | 330           | U |  |
| 2,6-dinitrotoluene        | 330           | U        | nitrobenzene               | 330           | Ù |  |
| di-n-octyl phthalate      | 330           | U        | n-nitroso-di-n-propylamine | 330           | U |  |
| 1,2-diphenylhydrazine(1)  | 330           | U        | n-nitrosodimethylamine     | 330           | U |  |
| fluoranthene              | 330           | U        | n-nitrosodiphenylamine(2)  | 330           | U |  |
| fluorene                  | 330           | U        | phenanthrene               | 330           | U |  |
| hexachlorobenzene         | 330           | Ų        | pyrene                     | 330           | U |  |
| hexachlorobutadiene       | 330           | U        | 1,2,4-trichlorobenzene     | .330          | U |  |
| hexachlorocyclopentadiene | 330           | U        |                            |               |   |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

<sup>(1) -</sup> Screened for as azobenzene

<sup>(2) -</sup> Detected as diphenylamine

Job Number: ITPK 51491

## ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in  $\mu g/kg$  (ppb)

Sample Matrix: Soil

Client Sample ID: Method Blank

Lab Sample ID: H0675

| Compound                   | Concen | tration | Compound              | Concer | ntration |
|----------------------------|--------|---------|-----------------------|--------|----------|
| 4-chloro-3-methylphenol    | 330    | U .     | 2-nitrophenol         | 330    | U        |
| 2-chlorophenol             | 330    | U .     | 4-nitrophenol         | 1,700  | U        |
| 2,4-dichlorophenol         | 330    | U       | pentachlorophenol     | 1,700  | U        |
| 2,4-dimethylphenol         | 330    | U       | phenol                | 330    | Ū        |
| 2,4-dinitrophenol          | 1,700  | Ŭ       | 2,4,6-trichlorophenol | 330    | U        |
| 2-methyl-4,6-dinitrophenol | 1,700  | U       |                       |        |          |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 06/

06/02/92

Date of Analysis:

06/08/92

J - Indicates an estimated value less than the detection limit.

Client Project ID: LE Carpenter

Job Number: ITPK 51491

## ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in  $\mu g/kg$  (ppb)

Sample Matrix: Soil

Client Sample ID: Method Blank

Lab Sample ID: H0675

| Tentative Identification (1)     | Concentration (2) |
|----------------------------------|-------------------|
| hexanedioic acid, dioctyl ester  | 3,800             |
| 2-pentanone, 4-hydroxy-4-methyl- | 6,500 A           |
| 2-pentanone, 5-(acetyloxy)-      | 640 À             |
| 2-heptanol, acetate              | 250               |
| cyclobutene, 2-propenylidene-    | 230               |

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

A - Suspected aldol condensation product.

Job Number: ITPK 51491

## SOIL SURROGATE PERCENT RECOVERY SUMMARY

## **SEMI-VOLATILE**

| Client Sample ID | Nitro-<br>Benzene-D5<br>(23-120%)* | 2-Fluoro-<br>Biphenyl<br>(30-115%)* | Terphenyl-<br>D14<br>(18-137%)* | Phenol-D5<br>(24-113%)* | 2-Fluoro-<br>Phenol<br>(25-121%)* | 2,4,6-<br>Tribromo-<br>Phenol<br>(19-122%)* |
|------------------|------------------------------------|-------------------------------------|---------------------------------|-------------------------|-----------------------------------|---|
| 1045-067-01      | 88                                 | 85                                  | 193 **                          | 98                      | 87                                | 40  |
| 1045-067-01 DL   | 80                                 | 80                                  | 124                             | 78                      | 64                                | 40  |
| 1045-067-02      | 90                                 | 88                                  | 105                             | 100                     | 85                                | 63  |
| 1045-067-03      | 88                                 | 83                                  | 96                              | 95                      | 75                                | 56  |
| 1045-067-04      | 95                                 | 88                                  | 125                             | 108                     | 91                                | 73  |
| 1045-067-04 DL   | D                                  | D                                   | D                               | 68                      | 66                                | D   |
| Method Blank     | 70                                 | 72                                  | 79                              | 71                      | 62                                | 62  |

<sup>\*</sup>Values in parenthesis represent QC limits.

<sup>\*\*</sup>Values outside of QC limits.

D - Surrogates diluted out DL - Dilution.

Client Project ID: LE Carpenter

Job Number: ITPK 51491

## TOTAL ORGANIC CARBON ANALYSIS

Results in mg/kg (ppm) dry weight

Sample Matrix: Soil

| Client Sample ID | t Sample ID Lab Sample ID |        |
|------------------|---------------------------|--------|
| Method Blank     | P3900                     | 1 U    |
| 1045-067-01      | K3228                     | 4,000  |
| 1045-067-02      | K3229                     | 11,000 |
| 1045-067-03      | K3230                     | 7,400  |
| 1045-067-04      | K3231                     | 11,000 |

Date of Analysis: 06/10/92

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

# Appendix F

Leachate Analyses

Job Number: ITAD 51019

## **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: 1045-055-1 Lab Sample ID: SS8501

|                          |               | •                         |          |       |
|--------------------------|---------------|---------------------------|----------|-------|
| Compound                 | Concentration | Compound                  | Concentr | ation |
| acrolein                 | 100 U         | 1,1-dichloroethene        | 50       | U     |
| acrylonitrile            | 100 U         | trans-1,2-dichloroethene  | 50       | U     |
| benzene                  | 110           | 1,2-dichloropropane       | 50       | U     |
| bromodichloromethane     | 50 U          | cis-1,3-dichloropropene   | 50       | U .   |
| bromoform                | <b>50</b> U   | trans-1,3-dichloropropene | 50       | U     |
| bromomethane             | 100 U         | ethyl benzene             | 1,700    |       |
| carbon tetrachloride     | 50 U          | methylene chloride        | 800      |       |
| chlorobenzene            | 50 U          | 1,1,2,2-tetrachloroethane | 50       | U     |
| chloroethane             | 100 U         | tetrachloroethene         | 50       | U     |
| 2-chloroethylvinyl ether | 100 U         | toluene                   | 320      |       |
| chloroform               | 30 J          | 1,1,1-trichloroethane     | 50       | U     |
| chloromethane            | 100 U         | 1,1,2-trichloroethane     | 50       | U     |
| dibromochloromethane     | 50 U          | trichloroethene           | 50       | U     |
| 1,1-dichloroethane       | <b>50</b> U   | trichlorofluoromethane    | 50       | U .   |
| 1,2-dichloroethane       | 50 Ü          | vinyl chloride            | 100      | U     |

Date of Analysis: 03/31/92

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Client Project ID: LE Carpenter

Job Number: ITAD 51019

## ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in  $\mu g/liter$  (ppb)

Sample Matrix: Water

Client Sample ID: 1045-055-1 Lab Sample ID: SS8501

Tentative Identification (1)

Concentration (2)

xylenes (total)

7,500

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- \* Quantitation based on continuing calibration.

Job Number: ITAD 51019

## **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: 1045-055-2 Lab Sample ID: SS8502

| Concentration | Compound   | Concent   | ration  |
|---------------|--|---|---|
| 10 U          | 1,1-dichloroethene                                       | 5   | U   |
| 10 U          | trans-1,2-dichloroethene                                 | .5  | U   |
| 120           | 1,2-dichloropropane                                      | 5   | U   |
| <b>2</b> J    | cis-1,3-dichloropropene                                  | 5   | U   |
| 5 Ü           | trans-1,3-dichloropropene                                | 5   | U   |
| 10 U          | ethyl benzene  | 71  |   |
| 5 U           | methylene chloride                                       | 320   | DB  |
| .5 U          | 1,1,2,2-tetrachloroethane                                | 5   | U   |
| 10 ; <b>U</b> | tetrachloroethene  | 5   | U   |
| 10 U          | toluene  | 310   | D   |
| 17            | 1,1,1-trichloroethane                                    | 5   | U   |
| 10 U          | 1,1,2-trichloroethane                                    | 5   | U   |
| 5 U           | trichloroethene  | 5   | U   |
| 5 U           | trichlorofluoromethane                                   | 5   | U   |
| 5 U           | vinyl chloride   | 10  | U   |
|               | 10 U 10 U 120 2 J 5 U 10 U 5 U 10 U 10 U 17 10 U 5 U 5 U | 10 U 1,1-dichloroethene  10 U trans-1,2-dichloroethene  120 1,2-dichloropropane  2 J cis-1,3-dichloropropene  5 U trans-1,3-dichloropropene  10 U ethyl benzene  5 U methylene chloride  5 U 1,1,2,2-tetrachloroethane  10 U toluene  17 1,1,1-trichloroethane  10 U 1,1,2-trichloroethane  5 U trichloroethene  10 U trichloroethane | 10       U       1,1-dichloroethene       5         10       U       trans-1,2-dichloroethene       5         120       1,2-dichloropropane       5         2       J       cis-1,3-dichloropropene       5         5       U       trans-1,3-dichloropropene       5         10       U       ethyl benzene       71         5       U       methylene chloride       320         5       U       1,1,2,2-tetrachloroethane       5         10       U       tetrachloroethene       5         10       U       toluene       310         17       1,1,1-trichloroethane       5         10       U       1,1,2-trichloroethane       5         5       U       trichloroethene       5         5       U       trichlorofluoromethane       5 |

Date of Analysis: 04/02/92

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

D - Compound analyzed at a secondary dilution factor.

Client Project ID: LE Carpenter

Job Number: ITAD 51019

## ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-055-2 Lab Sample ID: SS8502

| Tentative Identification (1) | Concentration (2) |
|------------------------------|-------------------|
| carbon disulfide             | 55 *              |
| acetone                      | 10 *              |
| xylenes (total)              | 840 *             |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- \* Quantitation based on continuing calibration.

Job Number: ITAD 51019

# **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: 1045-055-3 Lab Sample ID: SS8503

| Compound                 | Concentration | Compound                  | Concent  | ration |
|--------------------------|---------------|---------------------------|----------|--------|
| acrolein                 | 10 U          | 1,1-dichloroethene        | 5        | Ū      |
| acrylonitrile            | 10 Uʻ         | trans-1,2-dichloroethene  | 5        | U      |
| benzene                  | 48            | 1,2-dichloropropane       | . 5      | U      |
| bromodichloromethane     | 2 J           | cis-1,3-dichloropropene   | <b>5</b> | U      |
| bromoform                | 5 U           | trans-1,3-dichloropropene | 5        | U      |
| bromomethane             | 10 U          | ethyl benzene             | 4        | J      |
| carbon tetrachloride     | 5 U           | methylene chloride        | 6,600    | DB     |
| chlorobenzene            | 5 U           | 1,1,2,2-tetrachloroethane | 5        | U      |
| chloroethane             | 10 U          | tetrachloroethene         | 5        | Ü      |
| 2-chloroethylvinyl ether | 10 Ü          | toluene                   | 100      |        |
| chloroform               | 18            | 1,1,1-trichloroethane     | 5        | U      |
| chloromethane            | 10 U          | 1,1,2-trichloroethane     | 5        | U      |
| dibromochloromethane     | 5 U           | trichloroethene           | 5        | U      |
| 1,1-dichloroethane       | 5 U           | trichlorofluoromethane    | <b>5</b> | U      |
| 1,2-dichloroethane       | 5 U           | vinyl chloride            | 10       | U      |

Date of Analysis: 04/02/92

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

D - Compound analyzed at a secondary dilution factor.

Client Project ID: LE Carpenter

Job Number: ITAD 51019

## ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-055-3 Lab Sample ID: SS8503

| Tentative Identification (1)             | Concentra | Concentration (2) |  |
|--|-----------|-------------------|--|
| acetone                                  | 13        | *                 |  |
| carbon disulfide                         | 66        | *                 |  |
| xylenes (total)                          | 410       | *                 |  |
| 2-propenoic acid, 2-methyl-, methylester | 12        |                   |  |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- Quantitation based on continuing calibration.

Job Number: ITAD 51019

## **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: 1045-055-4 Lab Sample ID: SS8504

| Compound                 | Concentration  | Compound                  | Concentration |
|--------------------------|----------------|---------------------------|---------------|
| acrolein                 | 10 to <b>U</b> | 1,1-dichloroethene        | 5 U           |
| acrylonitrile            | 10 U           | trans-1,2-dichloroethene  | 5 U           |
| benzene                  | 10             | 1,2-dichloropropane       | 5 U           |
| bromodichloromethane     | 1 J            | cis-1,3-dichloropropene   | 5 U           |
| bromoform                | 5 U            | trans-1,3-dichloropropene | 5 U           |
| bromomethane             | 10 U           | ethyl benzene             | 5 U           |
| carbon tetrachloride     | 5 U            | methylene chloride        | 110 B         |
| chlorobenzene            | 5 U            | 1,1,2,2-tetrachloroethane | 5 U           |
| chloroethane             | 10 U           | tetrachloroethene         | 5 U           |
| 2-chloroethylvinyl ether | 10 Ü           | toluene                   | 2 J           |
| chloroform               | 13             | 1,1,1-trichloroethane     | 5 U           |
| chloromethane            | 10 Ü           | 1,1,2-trichloroethane     | 5 U           |
| dibromochloromethane     | 5 U            | trichloroethene           | 5 U           |
| 1,1-dichloroethane       | 5 U            | trichlorofluoromethane    | 5 U           |
| 1,2-dichloroethane       | 5 U            | vinyl chloride            | 10 U          |

Date of Analysis: 04/02/92

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

Client Project ID: LE Carpenter

Job Number: ITAD 51019

## ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-055-4 Lab Sample ID: SS8504

| <u>lentative identification (1)</u> | Concentration (2 | 1 |
|-------------------------------------|------------------|---|
| acetone                             | 3.7 *            |   |
| carbon disulfide                    | 9.0 *            |   |
| xylenes (total)                     | 64 *             |   |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- \* Quantitation based on continuing calibration.

Job Number: ITAD 51019

## **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: 1045-055-5 Lab Sample ID: SS8505

| Compound                 | Concentration | Compound                  | Concentration |
|--------------------------|---------------|---------------------------|---------------|
| acrolein                 | 10 U          | 1,1-dichloroethene        | 5 U           |
| acrylonitrile            | 10 U          | trans-1,2-dichloroethene  | 5 U           |
| benzene                  | 5 U           | 1,2-dichloropropane       | 5 Ü           |
| bromodichloromethane     | 8:            | cis-1,3-dichloropropene   | 5 U           |
| bromoform                | 5 U           | trans-1,3-dichloropropene | 5 U           |
| bromomethane             | 10 Ù          | ethyl benzene             | 5 U           |
| carbon tetrachloride     | 5 U           | methylene chloride        | 1 B           |
| chlorobenzene            | 5 U           | 1,1,2,2-tetrachloroethane | 5 U           |
| chloroethane             | 10 U          | tetrachloroethene         | 5 Ù           |
| 2-chloroethylvinyl ether | 10 U          | toluene                   | 5 U           |
| chloroform               | 67            | 1,1,1-trichloroethane     | 5 U           |
| chloromethane            | 10 U          | 1,1,2-trichloroethane     | 5 U           |
| dibromochloromethane     | 5 U           | trichloroethene           | 5 U           |
| 1,1-dichloroethane       | 5 U           | trichlorofluoromethane    | 5 U           |
| 1,2-dichloroethane       | 5 U           | vinyl chloride            | 10 U          |

Date of Analysis: 04/02/92

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

I - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

Client Project ID: LE Carpenter

Job Number: ITAD 51019

## ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-055-5 Lab Sample ID: SS8505

Tentative Identification (1)

Concentration (2)

none detected

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITAD 51019

## **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: Method Blank 1 Lab Sample ID: VB03313

| Compound                 | Concentration | Compound                  | Concentrat | ion |
|--------------------------|---------------|---------------------------|------------|-----|
| acrolein                 | 10 U          | 1,1-dichloroethene        | 5 t        | Ĵ   |
| acrylonitrile            | 10 U          | trans-1,2-dichloroethene  | 5 T        | J   |
| benzene                  | 5 U           | 1,2-dichloropropane       | 5 t        | J   |
| bromodichloromethane     | 5 U           | cis-1,3-dichloropropene   | 5 T        | J   |
| bromoform                | 5 U           | trans-1,3-dichloropropene | 5 T        | J   |
| bromomethane             | 10 U          | ethyl benzene             | 5 T        | J   |
| carbon tetrachloride     | 5 U           | methylene chloride        | 1 J        | Ī   |
| chlorobenzene            | 5 U           | 1,1,2,2-tetrachloroethane | 5 T        | J   |
| chloroethane             | 10 U          | tetrachloroethene         | 5 T        | J   |
| 2-chloroethylvinyl ether | 10 U          | toluene                   | <i>5</i> t | J,  |
| chloroform               | 5 U           | 1,1,1-trichloroethane     | 5 T        | J   |
| chloromethane            | 10 U          | 1,1,2-trichloroethane     | <i>5</i> t | J   |
| dibromochloromethane     | 5 U           | trichloroethene           | 5 T        | J   |
| 1,1-dichloroethane       | 5 ⊨ <b>U</b>  | trichlorofluoromethane    | 5 t        | Ĵ   |
| 1,2-dichloroethane       | 5 U           | vinyl chloride            | 10 U       | Ţ   |

Date of Analysis: 03/31/92

This method blank applies to the following sample: 1045-055-1.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Indicates an estimated value less than the detection limit.

Client Project ID: LE Carpenter

Job Number: ITAD 51019

## ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 1 Lab Sample ID: VB03313

Tentative Identification (1)

Concentration (2)

xylenes (total)

1.0

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- \* Quantitation based on continuing calibration.

Job Number: ITAD 51019

# **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: Method Blank 2

Lab Sample ID: EB0402

| Compound                 | Concent | ration   | Compound                  | Concent  | ration |
|--------------------------|---------|----------|---------------------------|----------|--------|
| acrolein                 | 10      | U        | 1,1-dichloroethene        | 5        | U      |
| acrylonitrile            | 10      | U ·      | trans-1,2-dichloroethene  | 5        | U      |
| benzene                  | 5       | U        | 1,2-dichloropropane       | 5        | U      |
| bromodichloromethane     | 5       | U        | cis-1,3-dichloropropene   | <b>5</b> | U      |
| bromoform                | 5       | U        | trans-1,3-dichloropropene | 5        | U      |
| bromomethane             | 10      | <b>U</b> | ethyl benzene             | Ś        | U      |
| carbon tetrachloride     | 5       | U .      | methylene chloride        | 1        | J      |
| chlorobenzene            | 5       | U        | 1,1,2,2-tetrachloroethane | 5        | U      |
| chloroethane             | 10      | U        | tetrachloroethene         | 5        | U      |
| 2-chloroethylvinyl ether | 10      | ់ប ់     | toluene                   | 5        | U      |
| chloroform               | 5       | Ù        | 1,1,1-trichloroethane     | 5        | U      |
| chloromethane            | 10      | U        | 1,1,2-trichloroethane     | 5        | U      |
| dibromochloromethane     | 5       | ប        | trichloroethene           | 5        | Ų.     |
| 1,1-dichloroethane       | 5       | U        | trichlorofluoromethane    | . 5      | U      |
| 1,2-dichloroethane       | 5       | U        | vinyl chloride            | 10       | U      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 04/02/92

This method blank applies to the following samples: 1045-055-1 DL, 1045-055-2, 1045-055-2 DL, 1045-055-3, 1045-055-3 DL 1045-055-4 and 1045-055-5.

DL - Dilution

J - Indicates an estimated value less than the detection limit.

Client Project ID: LE Carpenter

Job Number: ITAD 51019

## ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 2

Lab Sample ID: EB0402

Tentative Identification (1)

Concentration (2)

none detected

Remarks:

(1) Identification is based on computer search of the NIST Library.

(2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITAD 51019

# WATER SURROGATE PERCENT RECOVERY SUMMARY

|                  |                       | VOLATILE          | -                                   |
|------------------|-----------------------|-------------------|-------------------------------------|
| Client Sample ID | Toluene-D8 (88-110%)* | BFB<br>(86-115%)* | 1,2 Dichloroethane-D4<br>(76-114%)* |
| 1045-055-1       | 96                    | 106               | 97                                  |
| Method Blank 1   | <b>100</b>            | 101               | 97                                  |
| 1045-055-1 DL    | 97                    | 96                | 89                                  |
| 1045-055-2       | 89                    | 96                | 86                                  |
| 1045-055-2 DL    | 100                   | 100               | 93                                  |
| 1045-055-3       | 101                   | 101               | 92                                  |
| 1045-055-3 DL    | 97                    | 94                | 86                                  |
| 1045-055-4       | 98                    | 93                | 83                                  |
| 1045-055-5       | 100                   | 97                | 86                                  |
| Method Blank 2   | 101                   | 98                | 94                                  |

<sup>\*</sup>Values in parenthesis represent required QC limits.

DL - Dilution

Client Project ID: LE Carpenter

Job Number: ITAD 51019

## BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-055-1 Lab Sample ID: SS8506

| Compound                   | Concentration        | Compound                    | Concentration |
|----------------------------|----------------------|-----------------------------|---------------|
| acenaphthene               | 10 U                 | bis(2-chloroisopropyl)ether | . 10 U        |
| acenaphthylene             | 10 U                 | bis(2-ethylhexyl)phthalate  | 10 Ü          |
| anthracene                 | 10 U                 | 4-bromophenyl phenyl ether  | 10 U          |
| benzidine                  | 50 U                 | 2-chloronaphthalene         | 10 U          |
| benzo(a)anthracene         | 10 U                 | 4-chlorophenyl phenyl ether | 10 U          |
| benzo(b)fluoranthene       | 10 U                 | chrysene                    | 10 Ü          |
| benzo(k)fluoranthene       | 10 U                 | dibenz(a,h)anthracene       | 10 U          |
| benzo(a)pyrene             | 10 U                 | di-n-butylphthalate         | 20            |
| benzo(g,h,i)perylene       | 10 U                 | 1,2-dichlorobenzene         | 10 U          |
| butylbenzylphthalate       | 10 U                 | 1,3-dichlorobenzene         | 10 U          |
| bis(2-chloroethoxy)methane | <b>10</b> . <b>U</b> | 1,4-dichlorobenzene         | 10 U          |
| bis(2-chloroethyl)ether    | 10 U                 |                             |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Job Number: ITAD 51019

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-055-1 Lab Sample ID: SS8506

| Compound                  | Concentration | Compound                   | Concentration |
|---------------------------|---------------|----------------------------|---------------|
| 3,3'dichlorobenzidine     | <b>20</b> U   | hexachloroethane           | 10 U          |
| diethylphthalate          | 86            | indeno(1,2,3-cd)pyrene     | 10 U          |
| dimethyl phthalate        | 26            | isophorone                 | 10 U          |
| 2,4-dinitrotoluene        | 10 U          | naphthalene                | 3 J           |
| 2,6-dinitrotoluene        | 10 U          | nitrobenzene               | 10 U          |
| di-n-octyl phthalate      | 10 U          | n-nitroso-di-n-propylamine | 10 U          |
| 1,2-diphenylhydrazine(1)  | 10 U          | n-nitrosodimethylamine     | 10 U          |
| fluoranthene              | 10 U          | n-nitrosodiphenylamine(2)  | 10 U          |
| fluorene                  | 10 U          | phenanthrene               | 10 U          |
| hexachlorobenzene         | 10 U          | pyrene                     | 10 U          |
| hexachlorobutadiene       | 10 U          | 1,2,4-trichlorobenzene     | 10 U          |
| hexachlorocyclopentadiene | 10 Ü          | •                          |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

<sup>(1) -</sup> Screened for as azobenzene

<sup>(2) -</sup> Detected as diphenylamine

Client Project ID: LE Carpenter

Job Number: ITAD 51019

## ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in  $\mu g$ /liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-055-1 Lab Sample ID: SS8506

| Tentative Identification (1)                          | Concentration (2) |
|---|-------------------|
| benzene, (1-methylethyl)-                             | 30                |
| ethanol, 2-[(phenylmethyl)amino]-                     | 11                |
| ethyl methyl benzene                                  | 43 Y              |
| trimethyl benzene                                     | 21 Y              |
| trimethyl benzene                                     | 49 Y              |
| trimethyl benzene                                     | 19 Y              |
| 1,2-benzisothiazole                                   | 17                |
| unknown   | 8.8               |
| 1,3,2-dioxaphosphorinane, 5,5-dimethyl-2-phenoxy-,2-o | 27                |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Y - Indistinguishable isomer in tentatively identified compounds.

Job Number: ITAD 51019

## BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in  $\mu g$ /liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-055-2 Lab Sample ID: SS8507

| Compound                   | Concentration | Compound                    | Concentration |
|----------------------------|---------------|-----------------------------|---------------|
| acenaphthene               | 10 U          | bis(2-chloroisopropyl)ether | 10 U          |
| acenaphthylene             | 10 U          | bis(2-ethylhexyl)phthalate  | 10 U          |
| anthracene                 | 10 U          | 4-bromophenyl phenyl ether  | 10 U          |
| benzidine                  | 50 U          | 2-chloronaphthalene         | 10 U          |
| benzo(a)anthracene         | 10 U          | 4-chlorophenyl phenyl ether | 10 U          |
| benzo(b)fluoranthene       | 10 U          | chrysene                    | 10 U          |
| benzo(k)fluoranthene       | 10 U.         | dibenz(a,h)anthracene       | 10 U          |
| benzo(a)pyrene             | 10 Ü          | di-n-butylphthalate         | 10 U          |
| benzo(g,h,i)perylene       | 10 U          | 1,2-dichlorobenzene         | 10 U          |
| butylbenzylphthalate       | 10 U          | 1,3-dichlorobenzene         | 10 U          |
| bis(2-chloroethoxy)methane | 10 U          | 1,4-dichlorobenzene         | 10 U          |
| bis(2-chloroethyl)ether    | 10 U          |                             |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Job Number: ITAD 51019

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-055-2 Lab Sample ID: SS8507

| Compound                  | Concentration | Compound                   | <b>Concentration</b> |
|---------------------------|---------------|----------------------------|----------------------|
| 3,3'dichlorobenzidine     | 20 U          | hexachloroethane           | 10 U                 |
| diethylphthalate          | 11            | indeno(1,2,3-cd)pyrene     | 10 U                 |
| dimethyl phthalate        | 10 U          | isophorone                 | 10 U                 |
| 2,4-dinitrotoluene        | 10 U          | naphthalene                | 10 U                 |
| 2,6-dinitrotoluene        | 10 U          | nitrobenzene               | 10 U                 |
| di-n-octyl phthalate      | 10 U          | n-nitroso-di-n-propylamine | 10 U                 |
| 1,2-diphenylhydrazine(1)  | 10 U          | n-nitrosodimethylamine     | 10 U                 |
| fluoranthene              | 10 U          | n-nitrosodiphenylamine(2)  | 10 U                 |
| fluorene                  | 10 U          | phenanthrene               | 10 U                 |
| hexachlorobenzene         | 10 U          | pyrene                     | 10 U                 |
| hexachlorobutadiene       | 10 U          | 1,2,4-trichlorobenzene     | 10 U                 |
| hexachlorocyclopentadiene | 10 U          |                            |                      |

- U Compound was analyzed for but not detected. The number is the detection limit for the sample.
- Indicates an estimated value less than the detection limit.
- (1) Screened for as azobenzene
- (2) Detected as diphenylamine

Job Number: ITAD 51019

Client Project ID: LE Carpenter

## ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-055-2 Lab Sample ID: SS8507

| Tentative Identification (1)     | Concentration (2) |
|----------------------------------|-------------------|
| 2-pentanone, 4-hydroxy-4-methyl- | 13 AB             |
| 3-undecene, 6-methyl-, (E)-      | 20                |
| 1,2-benzisothiazole              | 15                |
| cyclohexane, isothiocyanato-     | . 11              |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- B Analyte was found in the blank as well as the sample.
- A Suspected aldol condensation product.

Job Number: ITAD 51019

## BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in  $\mu g$ /liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-055-3 Lab Sample ID: SS8508

| Compound                   | Concentration | Compound                    | Concentration |
|----------------------------|---------------|-----------------------------|---------------|
| acenaphthene               | 10 Ü          | bis(2-chloroisopropyl)ether | 10 U          |
| acenaphthylene             | 10 U          | bis(2-ethylhexyl)phthalate  | 10 U          |
| anthracene                 | 10 U          | 4-bromophenyl phenyl ether  | 10 U          |
| benzidine                  | <b>50 U</b>   | 2-chloronaphthalene         | 10 U          |
| benzo(a)anthracene         | 10 U          | 4-chlorophenyl phenyl ether | 10 U          |
| benzo(b)fluoranthene       | 10 U          | chrysene                    | 10 U          |
| benzo(k)fluoranthene       | 10 U          | dibenz(a,h)anthracene       | 10 U          |
| benzo(a)pyrene             | 10 U          | di-n-butylphthalate         | 10 U          |
| benzo(g,h,i)perylene       | 10 U          | 1,2-dichlorobenzene         | 10 U          |
| butylbenzylphthalate       | 10 Ü          | 1,3-dichlorobenzene         | 10 U          |
| bis(2-chloroethoxy)methane | 10 U          | 1,4-dichlorobenzene         | 10 U          |
| bis(2-chloroethyl)ether    | 10 U          | •                           |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Job Number: ITAD 51019

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-055-3 Lab Sample ID: SS8508

| Compound                  | Concentra | <u>tion</u> | Compound                   | Concentra | <u>ttion</u> |
|---------------------------|-----------|-------------|----------------------------|-----------|--------------|
| 3,3'dichlorobenzidine     | 20        | U.          | hexachloroethane           | 10        | U            |
| diethylphthalate          | i 12      |             | indeno(1,2,3-cd)pyrene     | 10        | U            |
| dimethyl phthalate        | 10        | Ü           | isophorone                 | 10        | U            |
| 2,4-dinitrotoluene        | 10        | U           | naphthalene                | 10        | U            |
| 2,6-dinitrotoluene        | 10        | U           | nitrobenzene               | 10        | U            |
| di-n-octyl phthalate      | 10        | U           | n-nitroso-di-n-propylamine | 10        | U            |
| 1,2-diphenylhydrazine(1)  | 10        | U           | n-nitrosodimethylamine     | 10        | Ü            |
| fluoranthene              | 10        | <b>U</b>    | n-nitrosodiphenylamine(2)  | 10        | U            |
| fluorene                  | 10        | U           | phenanthrene               | 10        | U            |
| hexachlorobenzene         | 10        | U           | pyrene                     | 10        | U            |
| hexachlorobutadiene       | 10        | U           | 1,2,4-trichlorobenzene     | 10        | U            |
| hexachlorocyclopentadiene | 10        | U           |                            |           |              |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

<sup>(1) -</sup> Screened for as azobenzene

<sup>(2) -</sup> Detected as diphenylamine

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE

KNOXVILLE, TN

Job Number: ITAD 51019

# Client Project ID: LE Carpenter

## ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-055-3 Lab Sample ID: SS8508

| Tentative Identification (1) | Concentration (2) |
|------------------------------|-------------------|
| 3-undecene, 6-methyl-, (E)-  | 25                |
| 1,2-benzisothiazole          | 14                |
| cyclohexane, isothiocyanato- | 17                |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITAD 51019

## BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-055-4 Lab Sample ID: SS8509

| Compound                   | Concentration | Compound                    | Concentration |
|----------------------------|---------------|-----------------------------|---------------|
| acenaphthene               | 10 U          | bis(2-chloroisopropyl)ether | 10 U          |
| acenaphthylene             | 10 U          | bis(2-ethylhexyl)phthalate  | 10 U          |
| anthracene                 | 10 U          | 4-bromophenyl phenyl ether  | 10 U          |
| benzidine                  | 50 U          | 2-chloronaphthalene         | 10 U          |
| benzo(a)anthracene         | 10 U          | 4-chlorophenyl phenyl ether | 10 U          |
| benzo(b)fluoranthene       | 10 U          | chrysene                    | 10 U          |
| benzo(k)fluoranthene       | 10 U          | dibenz(a,h)anthracene       | 10 U          |
| benzo(a)pyrene             | 10 U          | di-n-butylphthalate         | 10 Ü          |
| benzo(g,h,i)perylene       | 10 U          | 1,2-dichlorobenzene         | 10 U          |
| butylbenzylphthalate       | 10 U          | 1,3-dichlorobenzene         | 10 U          |
| bis(2-chloroethoxy)methane | 10 U          | 1,4-dichlorobenzene         | 10 U          |
| bis(2-chloroethyl)ether    | 10 U          |                             |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Job Number: ITAD 51019

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-055-4 Lab Sample ID: SS8509

| Compound                  | Concentration | Compound                   | Concentration |
|---------------------------|---------------|----------------------------|---------------|
| 3,3'dichlorobenzidine     | <b>20</b> U   | hexachloroethane           | 10 U          |
| diethylphthalate          | 2 J           | indeno(1,2,3-cd)pyrene     | 10 ·U         |
| dimethyl phthalate        | 10 U          | isophorone                 | 10 U          |
| 2,4-dinitrotoluene        | 10 U          | naphthalene                | 10 U          |
| 2,6-dinitrotoluene        | 10 U          | nitrobenzene               | 10 U          |
| di-n-octyl phthalate      | 10 U          | n-nitroso-di-n-propylamine | 10 U          |
| 1,2-diphenylhydrazine(1)  | 10 U          | n-nitrosodimethylamine     | 10 U          |
| fluoranthene              | 10 U          | n-nitrosodiphenylamine(2)  | 10 U          |
| fluorene                  | 10 U          | phenanthrene               | 10 U          |
| hexachlorobenzene         | 10 U          | pyrene                     | 10 U          |
| hexachlorobutadiene       | 10 U          | 1,2,4-trichlorobenzene     | 10 U          |
| hexachlorocyclopentadiene | 10 U          |                            |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

<sup>(1) -</sup> Screened for as azobenzene

<sup>(2) -</sup> Detected as diphenylamine

Client Project ID: LE Carpenter

Job Number: ITAD 51019

## ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-055-4 Lab Sample ID: SS8509

| Tentative Identification (1)     | Concentration (2) |
|----------------------------------|-------------------|
| 2-pentanone, 4-hydroxy-4-methyl- | 8.2 AB            |
| 3-undecene, 6-methyl-, (E)-      | . 11              |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- B Analyte was found in the blank as well as the sample.
- A Suspected aldol condensation product.

Job Number: ITAD 51019

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-055-5 Lab Sample ID: SS8510

| Compound                   | Concentration | Compound                                | Concentration |
|----------------------------|---------------|---|---------------|
| acenaphthene               | 10 Ü          | bis(2-chloroisopropyl)ether             | 10 U          |
| acenaphthylene             | 10 U          | bis(2-ethylhexyl)phthalate              | 10 U          |
| anthracene                 | 10 U          | 4-bromophenyl phenyl ether              | 10 U          |
| benzidine                  | <b>50</b> U   | 2-chloronaphthalene                     | 10 U          |
| benzo(a)anthracene         | 10 U          | 4-chlorophenyl phenyl ether             | 10 U          |
| benzo(b)fluoranthene       | 10 U          | chrysene                                | 10 U          |
| benzo(k)fluoranthene       | 10 U          | dibenz(a,h)anthracene                   | 10 U          |
| benzo(a)pyrene             | 10 U          | di-n-butylphthalate                     | 10 U          |
| benzo(g,h,i)perylene       | 10 U          | 1,2-dichlorobenzene                     | 10 Ŭ          |
| butylbenzylphthalate       | 10 U          | 1,3-dichlorobenzene                     | 10 U          |
| bis(2-chloroethoxy)methane | 10 U          | 1,4-dichlorobenzene                     | 10 U          |
| bis(2-chloroethyl)ether    | 10 U          | 1 · 5 · · · · · · · · · · · · · · · · · |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Job Number: ITAD 51019

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-055-5 Lab Sample ID: SS8510

| Compound                  | <u>Concentratio</u> | n Compound                 | Concentra | ation |
|---------------------------|---------------------|----------------------------|-----------|-------|
| 3,3'dichlorobenzidine     | 20 U                | hexachloroethane           | 10        | U     |
| diethylphthalate          | 10 U                | indeno(1,2,3-cd)pyrene     | 10        | U     |
| dimethyl phthalate        | 10 U                | isophorone                 | 10        | U     |
| 2,4-dinitrotoluene        | 10 U                | naphthalene                | 10        | U     |
| 2,6-dinitrotoluene        | 10 Ü                | nitrobenzene               | 10        | U     |
| di-n-octyl phthalate      | 10 U                | n-nitroso-di-n-propylamine | 10        | U     |
| 1,2-diphenylhydrazine(1)  | 10 U                | n-nitrosodimethylamine     | 10        | U     |
| fluoranthene              | 10 U                | n-nitrosodiphenylamine(2)  | 10        | U     |
| fluorene                  | 10 U                | phenanthrene               | 10        | U     |
| hexachlorobenzene         | 10 U                | pyrene                     | 10        | U     |
| hexachlorobutadiene       | 10 U                | 1,2,4-trichlorobenzene     | 10        | U     |
| hexachlorocyclopentadiene | 10 U                |                            |           |       |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

<sup>(1) -</sup> Screened for as azobenzene

<sup>(2) -</sup> Detected as diphenylamine

Client Project ID: LE Carpenter

Job Number: ITAD 51019

## ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-055-5 Lab Sample ID: SS8510

Tentative Identification (1)

Concentration (2)

none detected

Remarks:

(1) Identification is based on computer search of the NIST Library.

(2) Concentration is based on a response factor of 1.00 relative to the internal standard.

KNOXVILLE, TN

Job Number: ITAD 51019

## BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 1

Lab Sample ID: BL0143

| Compound                   | <b>Concentration</b> | Compound                    | Concentration |
|----------------------------|----------------------|-----------------------------|---------------|
| acenaphthene               | 10 U                 | bis(2-chloroisopropyl)ether | 10 Ù          |
| acenaphthylene             | 10 U                 | bis(2-ethylhexyl)phthalate  | 3 J           |
| anthracene                 | 10 U                 | 4-bromophenyl phenyl ether  | 10 U          |
| benzidine                  | 50 U                 | 2-chloronaphthalene         | 10 U          |
| benzo(a)anthracene         | 10 U                 | 4-chlorophenyl phenyl ether | 10 U          |
| benzo(b)fluoranthene       | 10 U                 | chrysene                    | 10 U          |
| benzo(k)fluoranthene       | 10 U                 | dibenz(a,h)anthracene       | 10 U          |
| benzo(a)pyrene             | 10 U                 | di-n-butylphthalate         | 10 U          |
| benzo(g,h,i)perylene       | 10 U                 | 1,2-dichlorobenzene         | 10 U          |
| butylbenzylphthalate       | 10 U                 | 1,3-dichlorobenzene         | 10 U          |
| bis(2-chloroethoxy)methane | 10 U                 | 1,4-dichlorobenzene         | 10 U          |
| bis(2-chloroethyl)ether    | 10 U                 |                             |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 03/27/92 Date of Analysis: 03/31/92

This method blank applies to the following samples: 1045-055-1, 1045-055-2, 1045-055-3, 1045-055-4 and 1045-055-5.

J - Indicates an estimated value less than the detection limit.

Job Number: ITAD 51019

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 1

Lab Sample ID: BL0143

|   | Compound                  | Concentra | <u>ttion</u> | Compound                   | <u>Concentra</u> | ation |
|---|---------------------------|-----------|--------------|----------------------------|------------------|-------|
|   | 3,3'dichlorobenzidine     | 20        | Ŭ            | hexachloroethane           | 10               | U     |
| • | diethylphthalate          | 10        | U ,          | indeno(1,2,3-cd)pyrene     | 10               | Ū     |
|   | dimethyl phthalate        | 10        | U            | isophorone                 | 10               | U     |
|   | 2,4-dinitrotoluene        | 10        | U .          | naphthalene                | 10               | U     |
|   | 2,6-dinitrotoluene        | 10        | Ü            | nitrobenzene               | 10               | U     |
|   | di-n-octyl phthalate      | 10        | U            | n-nitroso-di-n-propylamine | 10               | U     |
|   | 1,2-diphenylhydrazine(1)  | 10        | ับ           | n-nitrosodimethylamine     | 10               | U     |
|   | fluoranthene              | 10        | U            | n-nitrosodiphenylamine(2)  | 10               | U     |
|   | fluorene                  | 10        | Ú            | phenanthrene               | 10               | U     |
|   | hexachlorobenzene         | , 10      | U            | pyrene                     | 10               | U     |
|   | hexachlorobutadiene       | 10        | Ü            | 1,2,4-trichlorobenzene     | 10               | U     |
|   | hexachlorocyclopentadiene | 10        | U            |                            |                  |       |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

<sup>(1) -</sup> Screened for as azobenzene

<sup>(2) -</sup> Detected as diphenylamine

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

Job Number: ITAD 51019

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 1

Lab Sample ID: BL0143

Tentative Identification (1)

2-pentanone, 4-hydroxy-4-methyl
unknown

Concentration (2)

A

60

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

A - Suspected aldol condensation product.

Job Number: ITAD 51019

# WATER SURROGATE PERCENT RECOVERY SUMMARY

## **SEMI-VOLATILE**

| Client Sample ID | Nitro-<br>Benzene-D5<br>(35-114%)* | 2-Fluoro-<br>Biphenyl<br>(43-116%)* | Terphenyl-<br>D14<br>(33-141%)* | Phenol-D5<br>(10-94%)* | 2-Fluoro-<br>Phenol<br>(21-100%)* | 2,4,6-<br>Tribromo-<br>Phenol<br>(10-123%)* |
|------------------|------------------------------------|-------------------------------------|---------------------------------|------------------------|-----------------------------------|---|
| 1045-055-1       | 87                                 | 93                                  | 66                              | NA                     | NA                                | NA .  |
| 1045-055-2       | 79                                 | 87                                  | 64                              | NA                     | NA                                | NA  |
| 1045-055-3       | 80                                 | 88                                  | 66                              | NA                     | NA                                | NA  |
| 1045-055-4       | 82                                 | 86                                  | 79                              | NA                     | NA                                | NA  |
| 1045-055-5       | 84                                 | 91                                  | 84                              | NA                     | NA                                | NA  |
| Method Blank 1   | 79                                 | 86                                  | 92                              | NA                     | NA                                | NA.   |

<sup>\*</sup>Values in parenthesis represent contract required QC limits.

NA - Not applicable

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

Job Number: ITAD 51019

## CHEMICAL OXYGEN DEMAND

Results in mg/liter (ppm)

Sample Matrix: Water

| Client Sample ID | Lab Sample ID | Result |  |
|------------------|---------------|--------|--|
| 1045-055-5       | SS8511        | 8.7    |  |
| Method Blank     | -             | *      |  |

Date of Analysis: 03/31/92

<sup>\* -</sup> A method blank is not applicable for this analysis.

Job Number: ITAD 51254

# **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: 1045-061-1 Lab Sample ID: K1180

| Compound                 | Concentration | Compound                  | Concent    | ration |
|--------------------------|---------------|---------------------------|------------|--------|
| acrolein                 | 10 U          | 1,1-dichloroethene        | 5          | U      |
| acrylonitrile            | 10 U          | trans-1,2-dichloroethene  | 5          | U      |
| benzene                  | 32            | 1,2-dichloropropane       | <b>5</b> . | U      |
| bromodichloromethane     | 3 J           | cis-1,3-dichloropropene   | 5          | U      |
| bromoform                | 5 U           | trans-1,3-dichloropropene | 5          | U      |
| bromomethane             | 10 U          | ethyl benzene             | 5          | U      |
| carbon tetrachloride     | <b>5 U</b>    | methylene chloride        | 1,500      | DB     |
| chlorobenzene            | <b>5</b> Ü    | 1,1,2,2-tetrachloroethane | 5          | U      |
| chloroethane             | 10 U          | tetrachloroethene         | 5          | Ù.     |
| 2-chloroethylvinyl ether | 10 U          | toluene                   | 40         | ,      |
| chloroform               | 40            | 1,1,1-trichloroethane     | 5          | Ú      |
| chloromethane            | 10 U          | 1,1,2-trichloroethane     | 5          | U      |
| dibromochloromethane     | 5 U           | trichloroethene           | 5          | U      |
| 1,1-dichloroethane       | 5 U           | trichlorofluoromethane    | 5          | U      |
| 1,2-dichloroethane       | 5 U           | vinyl chloride            | 10         | U      |

Date of Analysis: 04/27/92

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

D - Compound analyzed at a secondary dilution factor.

Client Project ID: LE Carpenter

Job Number: ITAD 51254

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-061-1 Lab Sample ID: K1180

| Tentative Identification (1) | Concentration (2) |   |  |
|------------------------------|-------------------|---|--|
| carbon disulfide             | 61                |   |  |
| 2-propenoic acid, 2-methyl-, | 25                |   |  |
| benzene, dimethyl-           | 2,400             | Y |  |
| benzene, dimethyl-           | 430               | Y |  |
| benzene, 1-ethyl-methyl-     | 45                | Y |  |

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- Y Indistinguishable isomer in tentatively identified compounds.

Job Number: ITAD 51254

## **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu$ g/liter (ppb) Sample Matrix: Water

Client Sample ID: 1045-061-2 Lab Sample ID: K1181

|                          | - I              | 1                         |               |
|--------------------------|------------------|---------------------------|---------------|
| Compound                 | Concentration    | Compound                  | Concentration |
| acrolein                 | 10 U             | 1,1-dichloroethene        | 5 U           |
| acrylonitrile            | 10 U             | trans-1,2-dichloroethene  | 5 U           |
| benzene                  | <b>4 J</b>       | 1,2-dichloropropane       | 5 U           |
| bromodichloromethane     | 3 <sup>1</sup> J | cis-1,3-dichloropropene   | 5 U           |
| bromoform                | 5 U              | trans-1,3-dichloropropene | 5 U           |
| bromomethane             | 10 U             | ethyl benzene             | 5 U           |
| carbon tetrachloride     | 5 U              | methylene chloride        | 570 DB        |
| chlorobenzene            | 5 U              | 1,1,2,2-tetrachloroethane | 5 U           |
| chloroethane             | 10 U             | tetrachloroethene         | 5 U           |
| 2-chloroethylvinyl ether | 10 U             | toluene                   | 5             |
| chloroform               | 27 ·             | 1,1,1-trichloroethane     | 5 U           |
| chloromethane            | 10 U             | 1,1,2-trichloroethane     | 5 U           |
| dibromochloromethane     | 5 U              | trichloroethene           | 5 U           |
| 1,1-dichloroethane       | 5 U              | trichlorofluoromethane    | 5 U           |
| 1,2-dichloroethane       | 5 U              | vinyl chloride            | 10 U          |

Date of Analysis: 04/27/92

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

D - Compound analyzed at a secondary dilution factor.

Client Project ID: LE Carpenter

Job Number: ITAD 51254

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-061-2 Lab Sample ID: K1181

| Tentative Identification (1) | Concentration (2) |  |  |
|------------------------------|-------------------|--|--|
| carbon disulfide             | 1.3               |  |  |
| 2-propenoic acid, 2-methyl-, | 5.0               |  |  |
| benzene, dimethyl-           | 4.2 Y             |  |  |
| benzene, dimethyl-           | 33 Y              |  |  |
| benzene, dimethyl-           | 25 Y              |  |  |

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Y - Indistinguishable isomer in tentatively identified compounds.

Job Number: ITAD 51254

## **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: 1045-061-3 Lab Sample ID: K1182

| Compound                 | Concentration | Compound                  | Concent | ration     |
|--------------------------|---------------|---------------------------|---------|------------|
| acrolein                 | 10 U          | 1,1-dichloroethene        | 5       | Ù          |
| acrylonitrile            | 10 U          | trans-1,2-dichloroethene  | 5       | U          |
| benzene                  | 3 J           | 1,2-dichloropropane       | 5       | U          |
| bromodichloromethane     | 4 Ј           | cis-1,3-dichloropropene   | .5      | U          |
| bromoform                | 5 U           | trans-1,3-dichloropropene | 5       | U          |
| bromomethane             | 10 U          | ethyl benzene             | 5       | U          |
| carbon tetrachloride     | 5 U           | methylene chloride        | 8,100   | DB         |
| chlorobenzene            | 5 U           | 1,1,2,2-tetrachloroethane | 5       | U          |
| chloroethane             | 10 U          | tetrachloroethene         | 5       | Ù          |
| 2-chloroethylvinyl ether | 10 U          | toluene                   | 3       | <b>J</b> . |
| chloroform               | 39            | 1,1,1-trichloroethane     | 5       | U          |
| chloromethane            | 10 U          | 1,1,2-trichloroethane     | 5       | Ü          |
| dibromochloromethane     | 5 U           | trichloroethene           | 5       | U          |
| 1,1-dichloroethane       | 5 U           | trichlorofluoromethane    | 5       | U          |
| 1,2-dichloroethane       | 5 □ <b>U</b>  | vinyl chloride            | 10      | Ù          |

Date of Analysis: 04/27/92

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

D - Compound analyzed at a secondary dilution factor.

Client Project ID: LE Carpenter

Job Number: ITAD 51254

## ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-061-3 Lab Sample ID: K1182

| Tentative Identification (1) | Concentration (2) |
|------------------------------|-------------------|
| benzene, dimethyl-           | 88 Y              |
| benzene, dimethyl-           | 18 Y              |

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Y - Indistinguishable isomer in tentatively identified compounds.

Job Number: ITAD 51254

## **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: 1045-061-4 Lab Sample ID: K1183

| Compound                 | Concentration | Compound                  | Concent | ration |
|--------------------------|---------------|---------------------------|---------|--------|
| acrolein                 | 10 U          | 1,1-dichloroethene        | 5       | U      |
| acrylonitrile            | 10 U          | trans-1,2-dichloroethene  | 5       | U      |
| benzene                  | 5 U           | 1,2-dichloropropane       | 5       | U      |
| bromodichloromethane     | 3 J           | cis-1,3-dichloropropene   | 5       | U      |
| bromoform                | .5 U          | trans-1,3-dichloropropene | 5       | Ù      |
| bromomethane             | 10 U          | ethyl benzene             | 5       | U      |
| carbon tetrachloride     | 5 U           | methylene chloride        | 64      | В      |
| chlorobenzene            | 5 U           | 1,1,2,2-tetrachloroethane | 5       | U      |
| chloroethane             | 10 U          | tetrachloroethene         | 5       | Ü      |
| 2-chloroethylvinyl ether | 10 U          | toluene                   | .5      | U.     |
| chloroform               | 43            | 1,1,1-trichloroethane     | 5       | U      |
| chloromethane            | 10 U          | 1,1,2-trichloroethane     | 5       | U      |
| dibromochloromethane     | 5 U           | trichloroethene           | 5       | U      |
| 1,1-dichloroethane       | 5 U           | trichlorofluoromethane    | 5       | U      |
| 1,2-dichloroethane       | 5 U           | vinyl chloride            | 10      | U      |

Date of Analysis: 04/27/92

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

Client Project ID: LE Carpenter

Job Number: ITAD 51254

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-061-4 Lab Sample ID: K1183

Tentative Identification (1)

Concentration (2)

none detected

Remarks:

(1) Identification is based on computer search of the NIST Library.

(2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITAD 51254

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g/liter$  (ppb) Sample Matrix: Water

Client Sample ID: Method Blank 1 Lab Sample ID: VB0427

| Compound                 | Concentration | Compound                  | Concentration |
|--------------------------|---------------|---------------------------|---------------|
| acrolein                 | - 10 U        | 1,1-dichloroethene        | 5 U           |
| acrylonitrile            | 10 U          | trans-1,2-dichloroethene  | 5 U           |
| benzene                  | 5 U           | 1,2-dichloropropane       | 5 U           |
| bromodichloromethane     | <b>5</b> Ù    | cis-1,3-dichloropropene   | 5 U           |
| bromoform                | 5 / U         | trans-1,3-dichloropropene | 5 U           |
| bromomethane             | 10 U          | ethyl benzene             | 5 U           |
| carbon tetrachloride     | 5 U           | methylene chloride        | 2 J           |
| chlorobenzene            | 5 U           | 1,1,2,2-tetrachloroethane | 5 U           |
| chloroethane             | 10 U          | tetrachloroethene         | 5 U           |
| 2-chloroethylvinyl ether | 10 U          | toluene                   | <b>5 U</b> .  |
| chloroform               | 5 U           | 1,1,1-trichloroethane     | 5 U           |
| chloromethane            | 10 U          | 1,1,2-trichloroethane     | 5 U           |
| dibromochloromethane     | 5 U           | trichloroethene           | 5 U           |
| 1,1-dichloroethane       | 5 U           | trichlorofluoromethane    | 5 U           |
| 1,2-dichloroethane       | 5 U.          | vinyl chloride            | 10 Ü          |

Date of Analysis: 04/27/92

This method blank applies to the following samples: 1045-061-1, 1045-061-2, 1045-061-3 and 1045-061-4.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Client Project ID: LE Carpenter

Job Number: ITAD 51254

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 1 Lab Sample ID: VB0427

Tentative Identification (1)

Concentration (2)

none detected

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITAD 51254

## **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: Method Blank 2

Lab Sample ID: VB0428

| Compound                 | Concentration        | Compound                  | Concent | ration   |
|--------------------------|----------------------|---------------------------|---------|----------|
| acrolein                 | 10 U                 | 1,1-dichloroethene        | 5       | U        |
| acrylonitrile            | 10 U                 | trans-1,2-dichloroethene  | 5       | U        |
| benzene                  | <b>5</b> ** <b>U</b> | 1,2-dichloropropane       | 5       | U        |
| bromodichloromethane     | 5 U                  | cis-1,3-dichloropropene   | 5       | U        |
| bromoform                | 5 U                  | trans-1,3-dichloropropene | 5       | U        |
| bromomethane             | 10 U                 | ethyl benzene             | 5       | Ù        |
| carbon tetrachloride     | 5 Ü                  | methylene chloride        | 2       | J.       |
| chlorobenzene            | 5 TU                 | 1,1,2,2-tetrachloroethane | 5       | U        |
| chloroethane             | 10 U                 | tetrachloroethene         | 5       | U        |
| 2-chloroethylvinyl ether | 10 U                 | toluene                   | 5       | <b>U</b> |
| chloroform               | 5 U                  | 1,1,1-trichloroethane     | 5       | U        |
| chloromethane            | 10 U                 | 1,1,2-trichloroethane     | 5       | U        |
| dibromochloromethane     | 5 U                  | trichloroethene           | 5       | Ų        |
| 1,1-dichloroethane       | 5 U                  | trichlorofluoromethane    | 5       | U        |
| 1,2-dichloroethane       | 5 U                  | vinyl chloride            | 10      | U        |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 04/28/92

This method blank applies to the following samples: 1045-061-1 DL, 1045-061-2 DL, and 1045-061-3 DL.

DL - Dilution

J - Indicates an estimated value less than the detection limit.

Client Project ID: LE Carpenter

Job Number: ITAD 51254

### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 2

Lab Sample ID: VB0428

Tentative Identification (1)

Concentration (2)

none detected

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITAD 51254

## WATER SURROGATE PERCENT RECOVERY SUMMARY

|                  | <u></u>               | VOLATILE          |                                     |
|------------------|-----------------------|-------------------|-------------------------------------|
| Client Sample ID | Toluene-D8 (88-110%)* | BFB<br>(86-115%)* | 1,2 Dichloroethane-D4<br>(76-114%)* |
| 1045-061-1       | 100                   | 117 **            | 100                                 |
| 1045-061-1 DL    | 108                   | 111               | 103                                 |
| 1045-061-2       | 99                    | 104               | 99                                  |
| 1045-061-2 DL    | 104                   | 104               | 96                                  |
| 1045-061-3       | 101                   | 106               | 99                                  |
| 1045-061-3 DL    | 102                   | 103               | 100                                 |
| 1045-061-4       | 99                    | 104               | 99                                  |
| Method Blank 1   | 108                   | 109               | 109                                 |
| Method Blank 2   | 96                    | 96                | 91                                  |

<sup>\*</sup>Values in parenthesis represent contract required QC limits.

<sup>\*\*</sup>Values outside of contract required QC limits.

Job Number: ITAD 51254

#### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-061-1 Lab Sample ID: K1184

| Compound                   | Concentration | Compound                    | Concentration |
|----------------------------|---------------|-----------------------------|---------------|
| acenaphthene               | 10 Ü          | bis(2-chloroisopropyl)ether | 10 U          |
| acenaphthylene             | 10 U          | bis(2-ethylhexyl)phthalate  | 10 U          |
| anthracene                 | 10 U          | 4-bromophenyl phenyl ether  | 10 U          |
| benzidine                  | 50 U          | 2-chloronaphthalene         | 10 U          |
| benzo(a)anthracene         | 10 U          | 4-chlorophenyl phenyl ether | 10 U          |
| benzo(b)fluoranthene       | 10 U          | chrysene                    | 10 U          |
| benzo(k)fluoranthene       | 10 U          | dibenz(a,h)anthracene       | 10 U          |
| benzo(a)pyrene             | 10 U          | di-n-butylphthalate         | 10 U          |
| benzo(g,h,i)perylene       | 10 U          | 1,2-dichlorobenzene         | 10 U          |
| butylbenzylphthalate       | 10 U          | 1,3-dichlorobenzene         | 10 U          |
| bis(2-chloroethoxy)methane | io U          | 1,4-dichlorobenzene         | 10 U          |
| bis(2-chloroethyl)ether    | 10 U          |                             |               |

Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction:

04/28/92

Date of Analysis:

05/05/92

Indicates an estimated value less than the detection limit.

Job Number: ITAD 51254

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-061-1 Lab Sample ID: K1184

| Compound                  | Concentration | Compound                   | <u>Concentration</u> |
|---------------------------|---------------|----------------------------|----------------------|
| 3,3'dichlorobenzidine     | 20 U          | hexachloroethane           | 10 U                 |
| diethylphthalate          | 10 U          | indeno(1,2,3-cd)pyrene     | 10 Ū                 |
| dimethyl phthalate        | 10 U          | isophorone                 | 10 U                 |
| 2,4-dinitrotoluene        | 10 U          | naphthalene                | 2 J                  |
| 2,6-dinitrotoluene        | 10 U          | nitrobenzene               | 10 U                 |
| di-n-octyl phthalate      | 10 U          | n-nitroso-di-n-propylamine | 10 U                 |
| 1,2-diphenylhydrazine(1)  | 10 U          | n-nitrosodimethylamine     | 10 U                 |
| fluoranthene              | 10 U          | n-nitrosodiphenylamine(2)  | 10 U                 |
| fluorene                  | 10 U          | phenanthrene               | 10 U                 |
| hexachlorobenzene         | 10 U          | pyrene                     | 10 U                 |
| hexachlorobutadiene       | 10 U          | 1,2,4-trichlorobenzene     | 10 U                 |
| hexachlorocyclopentadiene | 10 U          |                            |                      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

<sup>(1) -</sup> Screened for as azobenzene

<sup>(2) -</sup> Detected as diphenylamine

Job Number: ITAD 51254

#### ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-061-1 Lab Sample ID: K1184

| Compound                   | Concen | tration | Compound              | Concer | ntration |
|----------------------------|--------|---------|-----------------------|--------|----------|
| 4-chloro-3-methylphenol    | 10     | U       | 2-nitrophenol         | 10     | U        |
| 2-chlorophenol             | 10     | U       | 4-nitrophenol         | 50     | U        |
| 2,4-dichlorophenol         | 10     | U       | pentachlorophenol     | 50     | Ü        |
| 2,4-dimethylphenol         | 130    |         | phenol                | 2      | J        |
| 2,4-dinitrophenol          | 50     | U       | 2,4,6-trichlorophenol | 10     | U        |
| 2-methyl-4,6-dinitrophenol | 50     | U       |                       |        |          |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction:

04/28/92

Date of Analysis:

05/05/92

J - Indicates an estimated value less than the detection limit.

Job Number: ITAD 51254

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-061-1 Lab Sample ID: K1184

| Tentative Identification (1) Concent |      |
|--------------------------------------|------|
| benzene, ethyl-methyl-               | 45 Y |
| benzene, ethyl-methyl-               | 24 Y |
| benzenemethanol, 2-methyl-           | 17   |
| benzene, trimethyl-                  | 22 Y |
| benzothiazole                        | 15   |
| benzene, trimethyl-                  | 16 Y |
| cyclohexane, isothiocyanato-         | 9.4  |

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Y - Indistinguishable isomer in tentatively identified compounds.

Job Number: ITAD 51254

## BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in  $\mu g$ /liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-061-2 Lab Sample ID: K1185

| Compound                   | Concentrat | ion      | Compound                    | Concentr | ation |
|----------------------------|------------|----------|-----------------------------|----------|-------|
| acenaphthene               | 10 U       | J        | bis(2-chloroisopropyl)ether | 10       | U     |
| acenaphthylene             | 10 U       | J        | bis(2-ethylhexyl)phthalate  | 10       | U     |
| anthracene                 | 10 U       | J        | 4-bromophenyl phenyl ether  | 10       | U     |
| benzidine                  | 50 U       | <b>J</b> | 2-chloronaphthalene         | 10       | U     |
| benzo(a)anthracene         | 10 U       | Ţ        | 4-chlorophenyl phenyl ether | 10       | U     |
| benzo(b)fluoranthene       | 10 U       | J        | chrysene                    | 10       | U     |
| benzo(k)fluoranthene       | 10 Ü       | Ţ        | dibenz(a,h)anthracene       | 10       | U     |
| benzo(a)pyrene             | 10 U       | Ţ        | di-n-butylphthalate         | 10       | U     |
| benzo(g,h,i)perylene       | 10 U       | <b>,</b> | 1,2-dichlorobenzene         | 10       | U     |
| butylbenzylphthalate       | 10 U       | Γ.       | 1,3-dichlorobenzene         | 10       | U     |
| bis(2-chloroethoxy)methane | 10 U       | Ţ        | 1,4-dichlorobenzene         | 10       | U     |
| bis(2-chloroethyl)ether    | 10 U       | ī        |                             | •        |       |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction:

04/28/92

Date of Analysis:

05/05/92

J - Indicates an estimated value less than the detection limit.

Job Number: ITAD 51254

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-061-2 Lab Sample ID: K1185

| <u>Compound</u>           | Concentration | Compound                   | Concentration |
|---------------------------|---------------|----------------------------|---------------|
| 3,3'dichlorobenzidine     | <b>20</b> U   | hexachloroethane           | 10 U          |
| diethylphthalate          | 10 U          | indeno(1,2,3-cd)pyrene     | 10 · U        |
| dimethyl phthalate        | 10 U          | isophorone                 | 10 U          |
| 2,4-dinitrotoluene        | 10 U          | naphthalene                | 10 U          |
| 2,6-dinitrotoluene        | 10 Ü          | nitrobenzene               | 10 U          |
| di-n-octyl phthalate      | 10 U          | n-nitroso-di-n-propylamine | 10 U          |
| 1,2-diphenylhydrazine(1)  | 10 Ù          | n-nitrosodimethylamine     | 10 U          |
| fluoranthene              | 10 U          | n-nitrosodiphenylamine(2)  | 10 U          |
| fluorene                  | 10 U          | phenanthrene               | 10 U          |
| hexachlorobenzene         | 10 U          | pyrene                     | 10 U          |
| hexachlorobutadiene       | 10 U          | 1,2,4-trichlorobenzene     | 10 U          |
| hexachlorocyclopentadiene | 10 U          |                            |               |

- U Compound was analyzed for but not detected. The number is the detection limit for the sample.
- J Indicates an estimated value less than the detection limit.
- (1) Screened for as azobenzene
- (2) Detected as diphenylamine

Job Number: ITAD 51254

# ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-061-2 Lab Sample ID: K1185

| Compound                   | Concentration | on Compound           | Concentration |
|----------------------------|---------------|-----------------------|---------------|
| 4-chloro-3-methylphenol    | 10 U          | 2-nitrophenol         | 10 U          |
| 2-chlorophenol             | 10 U          | 4-nitrophenol         | 50 U          |
| 2,4-dichlorophenol         | 10 U          | pentachlorophenol     | 50 U          |
| 2,4-dimethylphenol         | 10 U          | phenol                | 10 U          |
| 2,4-dinitrophenol          | <b>50</b> U   | 2,4,6-trichlorophenol | 10 U          |
| 2-methyl-4,6-dinitrophenol | 50 U          |                       |               |

- Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Date of Extraction:

04/28/92

Date of Analysis:

05/05/92

Client Project ID: LE Carpenter

Job Number: ITAD 51254

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-061-2 Lab Sample ID: K1185

Tentative Identification (1)

Concentration (2)

none detected

Remarks:

(1) Identification is based on computer search of the NIST Library.

(2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITAD 51254

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in  $\mu g$ /liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-061-3 Lab Sample ID: K1186

| Compound                   | Concentration | Compound                    | Concentration |
|----------------------------|---------------|-----------------------------|---------------|
| acenaphthene               | 10 U          | bis(2-chloroisopropyl)ether | 10 U          |
| acenaphthylene             | 10 U          | bis(2-ethylhexyl)phthalate  | 10 U          |
| anthracene                 | 10 U          | 4-bromophenyl phenyl ether  | 10 U          |
| benzidine                  | 50 U          | 2-chloronaphthalene         | 10 U          |
| benzo(a)anthracene         | 10 Ü          | 4-chlorophenyl phenyl ether | 10 U          |
| benzo(b)fluoranthene       | 10 U          | chrysene                    | 10 U          |
| benzo(k)fluoranthene       | 10 U          | dibenz(a,h)anthracene       | 10 U          |
| benzo(a)pyrene             | .10 Ù         | di-n-butylphthalate         | 10 U          |
| benzo(g,h,i)perylene       | 10 Û          | 1,2-dichlorobenzene         | 10 U          |
| butylbenzylphthalate       | 10 U          | 1,3-dichlorobenzene         | 10 U          |
| bis(2-chloroethoxy)methane | 10 U          | 1,4-dichlorobenzene         | 10 U          |
| bis(2-chloroethyl)ether    | 10 U          |                             | •             |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Job Number: ITAD 51254

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-061-3 Lab Sample ID: K1186

| Compound                  | Concentration | Compound                   | <b>Concentration</b> |
|---------------------------|---------------|----------------------------|----------------------|
| 3,3'dichlorobenzidine     | 20 U          | hexachloroethane           | 10 U                 |
| diethylphthalate          | 10 U          | indeno(1,2,3-cd)pyrene     | 10 U                 |
| dimethyl phthalate        | 10 U          | isophorone                 | 10 U                 |
| 2,4-dinitrotoluene        | 10 U          | naphthalene                | 10 U                 |
| 2,6-dinitrotoluene        | 10 Ŭ          | nitrobenzene               | 10 U                 |
| di-n-octyl phthalate      | 10 U          | n-nitroso-di-n-propylamine | 10 Ù                 |
| 1,2-diphenylhydrazine(1)  | 10 U          | n-nitrosodimethylamine     | 10 U                 |
| fluoranthene              | 10 U          | n-nitrosodiphenylamine(2)  | 10 U                 |
| fluorene                  | 10 U          | phenanthrene               | 10 U                 |
| hexachlorobenzene         | 10 U          | pyrene                     | , 10 <b>n</b>        |
| hexachlorobutadiene       | 10 U          | 1,2,4-trichlorobenzene     | 10 Ù                 |
| hexachlorocyclopentadiene | 10 U          |                            |                      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

<sup>(1) -</sup> Screened for as azobenzene

<sup>(2) -</sup> Detected as diphenylamine

Job Number: ITAD 51254

## ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in  $\mu g$ /liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-061-3 Lab Sample ID: K1186

| Compound                   | Concent | ration | Compound              | Concent | ration |
|----------------------------|---------|--------|-----------------------|---------|--------|
| 4-chloro-3-methylphenol    | 10      | U      | 2-nitrophenol         | 10      | Ū      |
| 2-chlorophenol             | 10      | U      | 4-nitrophenol         | 50      | U      |
| 2,4-dichlorophenol         | 10      | U      | pentachlorophenol     | 50      | U      |
| 2,4-dimethylphenol         | 10      | U      | phenol                | 10      | U      |
| 2,4-dinitrophenol          | 50      | U      | 2,4,6-trichlorophenol | 10      | U      |
| 2-methyl-4,6-dinitrophenol | 50      | U      |                       |         |        |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

IT Corporation May 29, 1992 IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

Client Project ID: LE Carpenter

Job Number: ITAD 51254

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in  $\mu g$ /liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-061-3 Lab Sample ID: K1186

Tentative Identification (1)

Concentration (2)

benzothiazole

11

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITAD 51254

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in  $\mu g/\text{liter}$  (ppb)

Sample Matrix: Water

Client Sample ID: 1045-061-4 Lab Sample ID: K1187

| Compound                   | Concentration | Compound                    | Concentration |
|----------------------------|---------------|-----------------------------|---------------|
| acenaphthene               | 10 U          | bis(2-chloroisopropyl)ether | 10 U          |
| acenaphthylene             | 10 U          | bis(2-ethylhexyl)phthalate  | 10 U          |
| anthracene                 | 10 U          | 4-bromophenyl phenyl ether  | 10 U          |
| benzidine                  | <b>50</b> U   | 2-chloronaphthalene         | 10 U          |
| benzo(a)anthracene         | 10 U          | 4-chlorophenyl phenyl ether | 10 U          |
| benzo(b)fluoranthene       | 10 U          | chrysene                    | 10 U          |
| benzo(k)fluoranthene       | 10 U          | dibenz(a,h)anthracene       | 10 U          |
| benzo(a)pyrene             | 10 U          | di-n-butylphthalate         | 10 U          |
| benzo(g,h,i)perylene       | 10 U          | 1,2-dichlorobenzene         | 10 .U         |
| butylbenzylphthalate       | 10 U          | 1,3-dichlorobenzene         | 10 U          |
| bis(2-chloroethoxy)methane | 10 U          | 1,4-dichlorobenzene         | 10 U          |
| bis(2-chloroethyl)ether    | 10 U          |                             |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 0

04/28/92

Date of Analysis:

05/05/92

J - Indicates an estimated value less than the detection limit.

Job Number: ITAD 51254

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-061-4 Lab Sample ID: K1187

| Compound                  | Concentration          | Compound                   | Concentration |  |  |
|---------------------------|------------------------|----------------------------|---------------|--|--|
| 3,3'dichlorobenzidine     | <b>20</b> U            | hexachloroethane           | 10 U          |  |  |
| diethylphthalate          | 10 U                   | indeno(1,2,3-cd)pyrene     | 10 U          |  |  |
| dimethyl phthalate        | 10 U                   | isophorone                 | 10 U          |  |  |
| 2,4-dinitrotoluene        | 10 U                   | naphthalene                | 10 U          |  |  |
| 2,6-dinitrotoluene        | 10 U                   | nitrobenzene               | 10 U          |  |  |
| di-n-octyl phthalate      | , <b>10</b> , <b>U</b> | n-nitroso-di-n-propylamine | 10 U          |  |  |
| 1,2-diphenylhydrazine(1)  | 10 U                   | n-nitrosodimethylamine     | 10 U          |  |  |
| fluoranthene              | 10 \ U                 | n-nitrosodiphenylamine(2)  | 10 U          |  |  |
| fluorene                  | 10 Ü                   | phenanthrene               | 10 U          |  |  |
| hexachlorobenzene         | 10 U                   | pyrene                     | 10 U          |  |  |
| hexachlorobutadiene       | 10 U                   | 1,2,4-trichlorobenzene     | 10 U          |  |  |
| hexachlorocyclopentadiene | 10 U                   |                            |               |  |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

<sup>(1) -</sup> Screened for as azobenzene

<sup>(2) -</sup> Detected as diphenylamine

Job Number: ITAD 51254

#### ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-061-4 Lab Sample ID: K1187

| Compound                   | Concent   | ration | Compound              | Concen | Concentration |  |
|----------------------------|-----------|--------|-----------------------|--------|---------------|--|
| 4-chloro-3-methylphenol    | 10        | U      | 2-nitrophenol         | 10     | U             |  |
| 2-chlorophenol             | 10        | U      | 4-nitrophenol         | 50     | U             |  |
| 2,4-dichlorophenol         | 10        | U      | pentachlorophenol     | 50     | Ū             |  |
| 2,4-dimethylphenol         | 10        | U      | phenol                | 10     | U             |  |
| 2,4-dinitrophenol          | <b>50</b> | Ü      | 2,4,6-trichlorophenol | 10     | U             |  |
| 2-methyl-4,6-dinitrophenol | 50        | U      |                       |        |               |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction:

04/28/92

Date of Analysis:

05/05/92

J - Indicates an estimated value less than the detection limit.

Client Project ID: LE Carpenter

Job Number: ITAD 51254

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-061-4 Lab Sample ID: K1187

Tentative Identification (1)

Concentration (2)

none detected

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITAD 51254

## BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank Lab Sample ID: H0436

| Compound                   | Concentration |          | Compound                    | Concentration |   |  |
|----------------------------|---------------|----------|-----------------------------|---------------|---|--|
| acenaphthene               | 10            | U        | bis(2-chloroisopropyl)ether | 10            | U |  |
| acenaphthylene             | 10            | U I      | bis(2-ethylhexyl)phthalate  | 10            | U |  |
| anthracene                 | 10            | U .      | 4-bromophenyl phenyl ether  | 10            | Ü |  |
| benzidine                  | 50            | Ü        | 2-chloronaphthalene         | 10            | U |  |
| benzo(a)anthracene         | 10            | U        | 4-chlorophenyl phenyl ether | 10            | U |  |
| benzo(b)fluoranthene       | 10            | <b>U</b> | chrysene                    | 10            | U |  |
| benzo(k)fluoranthene       | 10            | U        | dibenz(a,h)anthracene       | 10            | U |  |
| benzo(a)pyrene             | 10            | U        | di-n-butylphthalate         | 10            | U |  |
| benzo(g,h,i)perylene       | 10            | U        | 1,2-dichlorobenzene         | 10            | U |  |
| butylbenzylphthalate       | 10            | U        | 1,3-dichlorobenzene         | 10            | U |  |
| bis(2-chloroethoxy)methane | 10            | U        | 1,4-dichlorobenzene         | 10            | U |  |
| bis(2-chloroethyl)ether    | 10            | U        |                             |               |   |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Job Number: ITAD 51254

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in  $\mu g$ /liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank Lab Sample ID: H0436

| Compound                  | Concentration     | Compound                   | <u>Concentration</u> |  |  |
|---------------------------|-------------------|----------------------------|----------------------|--|--|
| 3,3'dichlorobenzidine     | 20 U              | hexachloroethane           | 10 U                 |  |  |
| diethylphthalate          | 10 <sup>'</sup> U | indeno(1,2,3-cd)pyrene     | 10 U                 |  |  |
| dimethyl phthalate        | 10 U              | isophorone                 | 10 U                 |  |  |
| 2,4-dinitrotoluene        | 10 U              | naphthalene                | 10 U                 |  |  |
| 2,6-dinitrotoluene        | 10 U              | nitrobenzene               | 10 U                 |  |  |
| di-n-octyl phthalate      | 10 U              | n-nitroso-di-n-propylamine | 10 U                 |  |  |
| 1,2-diphenylhydrazine(1)  | 10 U              | n-nitrosodimethylamine     | 10 U                 |  |  |
| fluoranthene              | 10 U              | n-nitrosodiphenylamine(2)  | 10 U                 |  |  |
| fluorene                  | 10 U              | phenanthrene               | 10 U                 |  |  |
| hexachlorobenzene         | 10 U              | pyrene                     | 10 U                 |  |  |
| hexachlorobutadiene       | 10 U              | 1,2,4-trichlorobenzene     | 10 U                 |  |  |
| hexachlorocyclopentadiene | 10 U              |                            |                      |  |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

<sup>(1) -</sup> Screened for as azobenzene

<sup>(2) -</sup> Detected as diphenylamine

Job Number: ITAD 51254

## ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank

Lab Sample ID: H0436

| Compound                   | Concen | tration | Compound              | Concentration |   |
|----------------------------|--------|---------|-----------------------|---------------|---|
| 4-chloro-3-methylphenol    | 10     | U .     | 2-nitrophenol         | 10            | U |
| 2-chlorophenol             | 10     | U       | 4-nitrophenol         | 50            | U |
| 2,4-dichlorophenol         | 10     | U       | pentachlorophenol     | 50            | Ū |
| 2,4-dimethylphenol         | 10     | U       | phenol                | 10            | U |
| 2,4-dinitrophenol          | 50     | U       | 2,4,6-trichlorophenol | 10            | U |
| 2-methyl-4,6-dinitrophenol | 50     | U       |                       |               |   |

J - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction:

04/28/92

Date of Analysis:

05/05/92

J - Indicates an estimated value less than the detection limit.

Client Project ID: LE Carpenter

Job Number: ITAD 51254

### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank

Lab Sample ID: H0436

Tentative Identification (1)

Concentration (2)

2-pentanone, 4-hydroxy-4-methyl-

16 A

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

A - Suspected aldol condensation product.

Job Number: ITAD 51254

# WATER SURROGATE PERCENT RECOVERY SUMMARY

# **SEMI-VOLATILE**

| Client Sample ID | Nitro-<br>Benzene-D5<br>(35-114%)* | 2-Fluoro-<br>Biphenyl<br>(43-116%)* | Terphenyl-<br>D14<br>(33-141%)* | Phenol-D5<br>(10-94%)* | 2-Fluoro-<br>Phenol<br>(21-100%)* | 2,4,6-<br>Tribromo-<br>Phenol<br>(10-123%)* |
|------------------|------------------------------------|-------------------------------------|---------------------------------|------------------------|-----------------------------------|---|
| 1045-061-1       | 84                                 | 97                                  | 96                              | 34                     | <i>5</i> 3                        | 68  |
| 1045-061-2       | 86                                 | 101                                 | 106                             | 27                     | 40                                | 60  |
| 1045-061-3       | 93                                 | 92                                  | 95                              | 35                     | 56                                | 61  |
| 1045-061-4       | 108                                | 99                                  | 90                              | 32                     | 46                                | 59  |
| Method Blank     | 89                                 | 101                                 | 113                             | 31                     | 42                                | 58  |

<sup>\*</sup>Values in parenthesis represent contract required QC limits.

Job Number: ITAD 51254

# CHEMICAL OXYGEN DEMAND

Results in mg/liter (ppm)

Sample Matrix: Water

| Client Sample ID | Lab Sample ID |       | Result |
|------------------|---------------|-------|--------|
| Method Blank     | .:            | -     | *      |
| 1045-061-1       | :             | K1188 | 25     |
| 1045-061-2       |               | K1189 | 6.2    |
| 1045-061-3       |               | K1190 | 15     |
| 1045-061-4       |               | K1191 | 12     |

Date of Analysis: 04/28/92

<sup>\* -</sup> A method blank is not applicable for this analysis.

Job Number: ITAD 51315

# **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: 1045-063-1 Lab Sample ID: K1746

| Compound                 | Concentration | Compound                  | Concent | ration |
|--------------------------|---------------|---------------------------|---------|--------|
| acrolein                 | 1,000 U       | 1,1-dichloroethene        | 500     | U      |
| acrylonitrile            | 1,000 U       | trans-1,2-dichloroethene  | 500     | U      |
| benzene                  | <b>500</b> U  | 1,2-dichloropropane       | 500     | U      |
| bromodichloromethane     | <b>500</b> U  | cis-1,3-dichloropropene   | 500     | U      |
| bromoform                | 500 U         | trans-1,3-dichloropropene | 500     | U      |
| bromomethane             | 1,000 U       | ethyl benzene             | 1,800   |        |
| carbon tetrachloride     | <b>500</b> U  | methylene chloride        | 1,100   | В      |
| chlorobenzene            | <b>500</b> U  | 1,1,2,2-tetrachloroethane | .500    | U      |
| chloroethane             | 1,000 U       | tetrachloroethene         | 500     | U      |
| 2-chloroethylvinyl ether | 1,000 U       | toluene                   | 500     | U      |
| chloroform               | 500 U         | 1,1,1-trichloroethane     | 500     | U      |
| chloromethane            | 1,000 U       | 1,1,2-trichloroethane     | 500     | U      |
| dibromochloromethane     | 500 U         | trichloroethene           | 500     | U      |
| 1,1-dichloroethane       | <b>500</b> U  | trichlorofluoromethane    | 500     | U      |
| 1,2-dichloroethane       | 500 U         | vinyl chloride            | 1,000   | U      |

Date of Analysis: 05/08/92

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

Job Number: ITAD 51315

### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-063-1 Lab Sample ID: K1746

| Tentative Identification (1)               | Concentration (2) |
|--|-------------------|
| benzene, dimethyl-                         | 14,000 Y          |
| nonane                                     | 1,500             |
| benzene, 1,2-dimethyl                      | 1,400             |
| octane, dimethyl-                          | 1,200 Y           |
| nonane, 3-methyl-                          | 2,600             |
| cyclohexane, diethyl                       | 1,500             |
| octane, dimethyl-                          | 4,300 Y           |
| nonane, 3-methyl-                          | 2,600             |
| cyclopentane, 1-methyl-3-(2-methylpropyl)- | 1,700             |
| decane                                     | 4,700             |

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Y - Indistinguishable isomer in tentatively identified compounds.

Job Number: ITAD 51315

# **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: 1045-063-2 Lab Sample ID: K1747

| Compound                 | Concentration | Compound                  | Concentra | ation |
|--------------------------|---------------|---------------------------|-----------|-------|
| acrolein                 | <b>50</b> U   | 1,1-dichloroethene        | 20        | U     |
| acrylonitrile            | <i>5</i> 0 U  | trans-1,2-dichloroethene  | 20        | U     |
| benzene                  | 7 J           | 1,2-dichloropropane       | 20        | Ų     |
| bromodichloromethane     | 20 U          | cis-1,3-dichloropropene   | 20        | U     |
| bromoform                | 20 U          | trans-1,3-dichloropropene | 20        | U     |
| bromomethane             | <b>50</b> U   | ethyl benzene             | 20        | U     |
| carbon tetrachloride     | 20 U          | methylene chloride        | 620       | В     |
| chlorobenzene            | 20 U          | 1,1,2,2-tetrachloroethane | 20        | U     |
| chloroethane             | <b>50</b> U   | tetrachloroethene         | 20        | U     |
| 2-chloroethylvinyl ether | <b>50</b> U   | toluene                   | 18        | J     |
| chloroform               | 8 J           | 1,1,1-trichloroethane     | 20        | U .   |
| chloromethane            | 50 U          | 1,1,2-trichloroethane     | 20        | U     |
| dibromochloromethane     | 20 U          | trichloroethene           | 20        | Ü     |
| 1,1-dichloroethane       | 20 U          | trichlorofluoromethane    | 20        | U     |
| 1,2-dichloroethane       | 20 U          | vinyl chloride            | 50        | U     |

Date of Analysis: 05/07/92

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

Client Project ID: LE Carpenter

Job Number: ITAD 51315

### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-063-2 Lab Sample ID: K1747

| Tentative Identification (1) | Concentration (2) |
|------------------------------|-------------------|
| acetone                      | 6.3               |
| benzene, dimethyl-           | 53 Y              |
| benzene, 1,2-dimethyl-       | 7.0               |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- Y Indistinguishable isomer in tentatively identified compounds.

Job Number: ITAD 51315

# **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: 1045-063-3 Lab Sample ID: K1748

| Compound                 | Concentration | Compound                  | Concent | ration |
|--------------------------|---------------|---------------------------|---------|--------|
| acrolein                 | 250 U         | 1,1-dichloroethene        | 120     | U      |
| acrylonitrile            | 250 U         | trans-1,2-dichloroethene  | 120     | U      |
| benzene                  | 120 U         | 1,2-dichloropropane       | 120     | U      |
| bromodichloromethane     | 120 U         | cis-1,3-dichloropropene   | 120     | U      |
| bromoform                | 120 U         | trans-1,3-dichloropropene | 120     | U      |
| bromomethane             | 250 U         | ethyl benzene             | 120     | U      |
| carbon tetrachloride     | 120 U         | methylene chloride        | 3,800   | В      |
| chlorobenzene            | 120 U         | 1,1,2,2-tetrachloroethane | 120     | U      |
| chloroethane             | 250 U         | tetrachloroethene         | 120     | U      |
| 2-chloroethylvinyl ether | 250 U         | toluene                   | 120     | U      |
| chloroform               | 46 J          | 1,1,1-trichloroethane     | 120     | U      |
| chloromethane            | 250 U         | 1,1,2-trichloroethane     | 120     | U      |
| dibromochloromethane     | 120 U         | trichloroethene           | 120     | U      |
| 1,1-dichloroethane       | 120 U         | trichlorofluoromethane    | 120     | U      |
| 1,2-dichloroethane       | 1 <b>20</b> U | vinyl chloride            | 250     | U      |

Date of Analysis: 05/07/92

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

Client Project ID: LE Carpenter

Job Number: ITAD 51315

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in  $\mu g$ /liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-063-3 Lab Sample ID: K1748

Tentative Identification (1)

Concentration (2)

benzene, dimethyl-

42 Y

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- Y Indistinguishable isomer in tentatively identified compounds.

Job Number: ITAD 51315

# **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: 1045-063-4 Lab Sample ID: K1749

| Compound                 |    | Concentration | Compound                  | Concentration |
|--------------------------|----|---------------|---------------------------|---------------|
| acrolein                 | 1  | 100 Ù         | 1,1-dichloroethene        | <b>50</b> U   |
| acrylonitrile            | a. | 100 🖟 U       | trans-1,2-dichloroethene  | 50 U          |
| benzene                  |    | 50 U          | 1,2-dichloropropane       | <b>50</b> U   |
| bromodichloromethane     |    | <b>50</b> U   | cis-1,3-dichloropropene   | 50 U          |
| bromoform                |    | <b>50</b> U   | trans-1,3-dichloropropene | <b>50</b> U   |
| bromomethane             |    | 100 U         | ethyl benzene             | 50 Ü          |
| carbon tetrachloride     |    | 50 U          | methylene chloride        | 64 B          |
| chlorobenzene            |    | 50 U          | 1,1,2,2-tetrachloroethane | 50 U          |
| chloroethane             |    | 100 U         | tetrachloroethene         | <b>50</b> U   |
| 2-chloroethylvinyl ether |    | 100 U         | toluene                   | <b>50</b> U   |
| chloroform               |    | 37 J          | 1,1,1-trichloroethane     | 50 U          |
| chloromethane            |    | 100 U         | 1,1,2-trichloroethane     | 50 U          |
| dibromochloromethane     |    | <b>50</b> U   | trichloroethene           | 50 U          |
| 1,1-dichloroethane       |    | <b>50</b> U   | trichlorofluoromethane    | 50 U          |
| 1,2-dichloroethane       |    | 50 U          | vinyl chloride            | 100 U         |

Date of Analysis: 05/07/92

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

Client Project ID: LE Carpenter

Job Number: ITAD 51315

# ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in  $\mu g/liter$  (ppb)

Sample Matrix: Water

Client Sample ID: 1045-063-4 Lab Sample ID: K1749

| Tentative Identification (1) | Concentration (2) |  |  |
|------------------------------|-------------------|--|--|
| benzene, dimethyl-           | 15 Y              |  |  |
| benzene, 1,2-dimethyl-       | 12                |  |  |

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Y - Indistinguishable isomer in tentatively identified compounds.

Job Number: ITAD 51315

### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: Method Blank 1

Lab Sample ID: VB0507

| Compound                 | Concent    | ration      | Compound                  | Concent | ration     |
|--------------------------|------------|-------------|---------------------------|---------|------------|
| acrolein                 | 10         | U           | 1,1-dichloroethene        | 5       | U          |
| acrylonitrile            | 10         | U           | trans-1,2-dichloroethene  | 5       | U          |
| benzene                  | 5          | Ū           | 1,2-dichloropropane       | 5.      | U          |
| bromodichloromethane     | 5          | U           | cis-1,3-dichloropropene   | 5       | U          |
| bromoform                | 5          | Ú           | trans-1,3-dichloropropene | 5       | U          |
| bromomethane             | 10         | U           | ethyl benzene             | 5       | U          |
| carbon tetrachloride     | 5          | U           | methylene chloride        | 1       | j          |
| chlorobenzene            | 5          | U           | 1,1,2,2-tetrachloroethane | 5       | U          |
| chloroethane             | 10         | U           | tetrachloroethene         | 5       | U          |
| 2-chloroethylvinyl ether | 10         | U           | toluene                   | 5       | <b>U</b> , |
| chloroform               | 5          | U           | 1,1,1-trichloroethane     | 5       | U          |
| chloromethane            | 10         | υ           | 1,1,2-trichloroethane     | 5       | U          |
| dibromochloromethane     | 5          | U           | trichloroethene           | 5       | U          |
| 1,1-dichloroethane       | <b>5</b> , | U           | trichlorofluoromethane    | 5       | U          |
| 1,2-dichloroethane       | 5          | <b>"U</b> " | vinyl chloride            | 10      | Ù          |

Date of Analysis: 05/07/92

This method blank applies to the following samples: 1045-063-2, 1045-063-3 and 1045-063-4.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Client Project ID: LE Carpenter

Job Number: ITAD 51315

# ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 1 Lab Sample ID: VB0507

Tentative Identification (1)

Concentration (2)

none detected

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITAD 51315

# VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: Method Blank 2

Lab Sample ID: EB0508

| Compound                 | Concentration    | Compound                  | Concent | ration     |
|--------------------------|------------------|---------------------------|---------|------------|
| acrolein                 | 10 U             | 1,1-dichloroethene        | . 5     | U          |
| acrylonitrile            | 10 U             | trans-1,2-dichloroethene  | 5       | Ų          |
| benzene                  | <b>5</b> † Û     | 1,2-dichloropropane       | 5       | U          |
| bromodichloromethane     | 5 <sup>‡</sup> U | cis-1,3-dichloropropene   | 5       | U          |
| bromoform                | 5 U              | trans-1,3-dichloropropene | 5       | U          |
| bromomethane             | 10 U             | ethyl benzene             | 5       | U          |
| carbon tetrachloride     | 5 U              | methylene chloride        | 1       | J          |
| chlorobenzene            | 5 U              | 1,1,2,2-tetrachloroethane | . 5     | U          |
| chloroethane             | 10 Ü             | tetrachloroethene         | 5       | U          |
| 2-chloroethylvinyl ether | 10 U             | toluene                   | 5       | <b>U</b> . |
| chloroform               | 5 U              | 1,1,1-trichloroethane     | 5       | U          |
| chloromethane            | 10 U             | 1,1,2-trichloroethane     | 5       | U          |
| dibromochloromethane     | 5 U              | trichloroethene           | 5       | U          |
| 1,1-dichloroethane       | 5 U              | trichlorofluoromethane    | 5       | U          |
| 1,2-dichloroethane       | 5 U              | vinyl chloride            | 10      | U          |

Date of Analysis: 05/08/92

This method blank applies to the following sample: 1045-063-1.

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Client Project ID: LE Carpenter

Job Number: ITAD 51315

### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 2

Lab Sample ID: EB0508

Tentative Identification (1)

Concentration (2)

none detected

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITAD 51315

# WATER SURROGATE PERCENT RECOVERY SUMMARY

|                  |                       | VOLATILE          |                                     |  |
|------------------|-----------------------|-------------------|-------------------------------------|--|
| Client Sample ID | Toluene-D8 (88-110%)* | BFB<br>(86-115%)* | 1,2 Dichloroethane-D4<br>(76-114%)* |  |
| 1045-063-2       | 97                    | 93                | 88                                  |  |
| 1045-063-3       | 98                    | 93                | 91                                  |  |
| 1045-063-4       | 96                    | 91                | 91                                  |  |
| Method Blank 1   | 103                   | 97                | 97                                  |  |
| 1045-063-1       | 100                   | 104               | 93                                  |  |
| Method Blank 2   | 103                   | 102               | 97                                  |  |

<sup>\*</sup>Values in parenthesis represent contract required QC limits.

Job Number: ITAD 51315

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-063-1 Lab Sample ID: K1754

| Compound                   | Concentration |          | Compound                    | <b>Concentration</b> |    |  |
|----------------------------|---------------|----------|-----------------------------|----------------------|----|--|
| acenaphthene               | 500           | U        | bis(2-chloroisopropyl)ether | 500                  | U  |  |
| acenaphthylene             | 500           | U        | bis(2-ethylhexyl)phthalate  | 61,000               | D  |  |
| anthracene                 | 500           | U        | 4-bromophenyl phenyl ether  | 500                  | Ü  |  |
| benzidine                  | 2,500         | U        | 2-chloronaphthalene         | 500                  | U  |  |
| benzo(a)anthracene         | 500           | U        | 4-chlorophenyl phenyl ether | 500                  | U  |  |
| benzo(b)fluoranthene       | 500           | U        | chrysene                    | 500                  | Ū  |  |
| benzo(k)fluoranthene       | 500           | U        | dibenz(a,h)anthracene       | 500                  | U  |  |
| benzo(a)pyrene             | 500           | <b>U</b> | di-n-butylphthalate         | 500                  | U  |  |
| benzo(g,h,i)perylene       | 500           | Ü        | 1,2-dichlorobenzene         | 500                  | U  |  |
| butylbenzylphthalate       | 540           |          | 1,3-dichlorobenzene         | 500                  | U  |  |
| bis(2-chloroethoxy)methane | 500           | U        | 1,4-dichlorobenzene         | 500                  | U. |  |
| bis(2-chloroethyl)ether    | 500           | U        |                             |                      |    |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

D - Compound analyzed at a secondary dilution factor.

Job Number: ITAD 51315

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-063-1 Lab Sample ID: K1754

| Compound                  | Concentration |            | Compound                   | Concentration |   |  |
|---------------------------|---------------|------------|----------------------------|---------------|---|--|
| 3,3'dichlorobenzidine     | 1,000         | U          | hexachloroethane           | 500           | U |  |
| diethylphthalate          | 500           | U          | indeno(1,2,3-cd)pyrene     | 500           | U |  |
| dimethyl phthalate        | 500           | Ů          | isophorone                 | 500           | U |  |
| 2,4-dinitrotoluene        | 500           | Ü          | naphthalene                | 500           | U |  |
| 2,6-dinitrotoluene        | 500           | U          | nitrobenzene               | 500           | U |  |
| di-n-octyl phthalate      | 190           | J          | n-nitroso-di-n-propylamine | 500           | U |  |
| 1,2-diphenylhydrazine(1)  | 500           | Ü          | n-nitrosodimethylamine     | 500           | U |  |
| fluoranthene              | 500           | <b>. U</b> | n-nitrosodiphenylamine(2)  | 500           | Ú |  |
| fluorene                  | 500           | U          | phenanthrene               | 500           | U |  |
| hexachlorobenzene         | 500           | U          | pyrene                     | 500           | U |  |
| hexachlorobutadiene       | 500           | U          | 1,2,4-trichlorobenzene     | 500           | U |  |
| hexachlorocyclopentadiene | 500           | U          |                            |               |   |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

<sup>(1) -</sup> Screened for as azobenzene

<sup>(2) -</sup> Detected as diphenylamine

Job Number: ITAD 51315

# ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in  $\mu g$ /liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-063-1 Lab Sample ID: K1754

| Compound                   | Concentration |          | Compound              | Concentration |   |  |
|----------------------------|---------------|----------|-----------------------|---------------|---|--|
| 4-chloro-3-methylphenol    | 500           | U        | 2-nitrophenol         | 500           | U |  |
| 2-chlorophenol             | 500           | U        | 4-nitrophenol         | 2,500         | U |  |
| 2,4-dichlorophenol         | 500           | <b>U</b> | pentachlorophenol     | 2,500         | U |  |
| 2,4-dimethylphenol         | 500           | Ū        | phenol                | 500           | U |  |
| 2,4-dinitrophenol          | 2,500         | <b>u</b> | 2,4,6-trichlorophenol | 500           | U |  |
| 2-methyl-4,6-dinitrophenol | 2,500         | <b>U</b> |                       |               |   |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Job Number: ITAD 51315

# ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-063-1 Lab Sample ID: K1754

| Tentative Identification (1) | Concentration (2) |
|------------------------------|-------------------|
| unknown                      | 4,700             |
| unknown                      | 21,000            |
| ethanol, 2-(dodecyloxy)-     | 21,000            |
| 1-dodecanol                  | 13,000            |
| unknown                      | 34,000            |
| ethanol, 2-(dodecyloxy)-     | 17,000            |
| unknown                      | 3,500             |
| unknown                      | 3,100             |
| unknown                      | 2,000             |
| 1-hexadecanol                | 5,400             |
| unknown                      | 4,500             |
| ethanol, 2-(dodecyloxy)-     | 4,500             |
| ethanol, 2-(hexadecyloxy)-   | 3,900             |
| decane                       | 4,000             |
| unknown                      | 6,300             |
| octane, 2,5-dimethyl-        | 1,700             |
| unknown                      | 3,000             |
| 1-nonene, 4,6,8-trimethyl-   | 1,600             |
| nonane, 3-methyl-            | 1,400             |
| 1-decene, dimethyl-          | 1,300 Y           |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Y - Indistinguishable isomer in tentatively identified compounds.

Job Number: ITAD 51315

### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in  $\mu g/liter$  (ppb)

Sample Matrix: Water

Client Sample ID: 1045-063-2 Lab Sample ID: K1755R

| Compound                   | Concentration | 1   | Compound                    | Concentr | ation |
|----------------------------|---------------|-----|-----------------------------|----------|-------|
| acenaphthene               | 10 U          |     | bis(2-chloroisopropyl)ether | 10       | U     |
| acenaphthylene             | 10 U          |     | bis(2-ethylhexyl)phthalate  | 10       | U     |
| anthracene                 | 10 U          |     | 4-bromophenyl phenyl ether  | 10       | Ü     |
| benzidine                  | 50 U          | 1   | 2-chloronaphthalene         | 10       | U     |
| benzo(a)anthracene         | 10 U          |     | 4-chlorophenyl phenyl ether | 10       | U     |
| benzo(b)fluoranthene       | 10 U          |     | chrysene                    | 10       | U     |
| benzo(k)fluoranthene       | 10 U          |     | dibenz(a,h)anthracene       | 10       | U     |
| benzo(a)pyrene             | 10 U          |     | di-n-butylphthalate         | 10       | U     |
| benzo(g,h,i)perylene       | 10 U          | 1.1 | 1,2-dichlorobenzene         | 10       | U     |
| butylbenzylphthalate       | 10 U          |     | 1,3-dichlorobenzene         | 10       | U     |
| bis(2-chloroethoxy)methane | 10 U          |     | 1,4-dichlorobenzene         | 10       | U     |
| bis(2-chloroethyl)ether    | 10 U          |     |                             |          |       |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Job Number: ITAD 51315

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-063-2 Lab Sample ID: K1755R

|   | Compound                  | Concentration |                  | Compound                   | Concentration |    |  |  |  |
|---|---------------------------|---------------|------------------|----------------------------|---------------|----|--|--|--|
|   | 3,3'dichlorobenzidine     | 20            | U                | hexachloroethane           | 10            | U  |  |  |  |
| • | diethylphthalate          | 10            | U                | indeno(1,2,3-cd)pyrene     | 10            | U  |  |  |  |
|   | dimethyl phthalate        | 10            | U                | isophorone                 | 10            | U  |  |  |  |
|   | 2,4-dinitrotoluene        | 10            | U                | naphthalene                | 10            | U, |  |  |  |
|   | 2,6-dinitrotoluene        | 10            | Ü                | nitrobenzene               | 10            | U  |  |  |  |
|   | di-n-octyl phthalate      | 10            | U                | n-nitroso-di-n-propylamine | 10            | U  |  |  |  |
|   | 1,2-diphenylhydrazine(1)  | 10            | <b>U</b>         | n-nitrosodimethylamine     | 10            | Ù  |  |  |  |
|   | fluoranthene              | 10            | <b>U</b>         | n-nitrosodiphenylamine(2)  | 10            | U  |  |  |  |
|   | flüorene                  | 10            | Ù                | phenanthrene               | 10            | U  |  |  |  |
|   | hexachlorobenzene         | 10            | $\mathbf{U}_{i}$ | pyrene                     | 10            | U  |  |  |  |
|   | hexachlorobutadiene       | 10            | , <b>U</b>       | 1,2,4-trichlorobenzene     | 10            | Ù  |  |  |  |
|   | hexachlorocyclopentadiene | 10            | U                |                            |               |    |  |  |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

<sup>(1) -</sup> Screened for as azobenzene

<sup>(2) -</sup> Detected as diphenylamine

Job Number: ITAD 51315

### ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-063-2 Lab Sample ID: K1755R

| Compound                   | Concen | Concentration |   | Compound              | Concentr |   |
|----------------------------|--------|---------------|---|-----------------------|----------|---|
| 4-chloro-3-methylphenol    | 10     | U             |   | 2-nitrophenol         | 10       | U |
| 2-chlorophenol             | 10     | Ù             | í | 4-nitrophenol         | 50       | U |
| 2,4-dichlorophenol         | 10     | U             |   | pentachlorophenol     | 50       | Ŭ |
| 2,4-dimethylphenol         | 10     | U             |   | phenol                | 10       | U |
| 2,4-dinitrophenol          | 50     | Ū             |   | 2,4,6-trichlorophenol | 10       | U |
| 2-methyl-4,6-dinitrophenol | 50     | U             |   |                       |          |   |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Client Project ID: LE Carpenter

Job Number: ITAD 51315

### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-063-2 Lab Sample ID: K1755R

Tentative Identification (1)

2-pentanone, 4-hydroxy-4-methylbenzothiazole

Concentration (2)

14 B

8.4

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- B Analyte was found in the blank as well as the sample.

Job Number: ITAD 51315

### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-063-3 Lab Sample ID: K1756R

| Compound                   | Concentration | Compound                    | Concentration |  |  |
|----------------------------|---------------|-----------------------------|---------------|--|--|
| acenaphthene               | 10 U          | bis(2-chloroisopropyl)ether | 10 U          |  |  |
| acenaphthylene             | 10 U          | bis(2-ethylhexyl)phthalate  | 10 U          |  |  |
| anthracene                 | 10 U          | 4-bromophenyl phenyl ether  | 10 U          |  |  |
| benzidine                  | 50 U          | 2-chloronaphthalene         | 10 U          |  |  |
| benzo(a)anthracene         | 10 U          | 4-chlorophenyl phenyl ether | 10 U          |  |  |
| benzo(b)fluoranthene       | 10 Ŭ          | chrysene                    | 10 U          |  |  |
| benzo(k)fluoranthene       | 10 U          | dibenz(a,h)anthracene       | 10 U          |  |  |
| benzo(a)pyrene             | 10 U          | di-n-butylphthalate         | 10 U          |  |  |
| benzo(g,h,i)perylene       | 10 U          | 1,2-dichlorobenzene         | 10 U          |  |  |
| butylbenzylphthalate       | 10 U          | 1,3-dichlorobenzene         | 10 U          |  |  |
| bis(2-chloroethoxy)methane | 10 U          | 1,4-dichlorobenzene         | 10 U          |  |  |
| bis(2-chloroethyl)ether    | 10 U          | •                           |               |  |  |

<sup>-</sup> Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction:

05/08/92

Date of Analysis:

05/13/92

Indicates an estimated value less than the detection limit.

Job Number: ITAD 51315

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-063-3 Lab Sample ID: K1756R

| Compound                  | Concentration |    | Compound                   | Concentration |    |  |
|---------------------------|---------------|----|----------------------------|---------------|----|--|
| 3,3'dichlorobenzidine     | 20            | U. | hexachloroethane           | 10            | U  |  |
| diethylphthalate          | 10            | U  | indeno(1,2,3-cd)pyrene     | 10            | ·U |  |
| dimethyl phthalate        | 10            | U  | isophorone                 | 10            | U  |  |
| 2,4-dinitrotoluene        | 10            | U. | naphthalene                | 10            | U  |  |
| 2,6-dinitrotoluene        | 10            | U  | nitrobenzene               | 10            | U  |  |
| di-n-octyl phthalate      | 10            | U  | n-nitroso-di-n-propylamine | 10            | U  |  |
| 1,2-diphenylhydrazine(1)  | 10            | U  | n-nitrosodimethylamine     | 10            | U  |  |
| fluoranthene              | 10            | U  | n-nitrosodiphenylamine(2)  | 10            | U  |  |
| fluorene                  | 10            | Ų  | phenanthrene               | 10            | U  |  |
| hexachlorobenzene         | 10            | U  | pyrene                     | 10            | U  |  |
| hexachlorobutadiene       | 10            | U  | 1,2,4-trichlorobenzene     | 10            | U  |  |
| hexachlorocyclopentadiene | 10            | U  |                            |               |    |  |
|                           |               |    |                            |               |    |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Indicates an estimated value less than the detection limit.

<sup>(1) -</sup> Screened for as azobenzene

<sup>(2) -</sup> Detected as diphenylamine

Job Number: ITAD 51315

### ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-063-3 Lab Sample ID: K1756R

| Compound                   | Concen     | tration | Compound              | ! | Concentration |   |
|----------------------------|------------|---------|-----------------------|---|---------------|---|
| 4-chloro-3-methylphenol    | 10         | U       | 2-nitrophenol         |   | 10            | U |
| 2-chlorophenol             | 10         | U       | 4-nitrophenol         |   | 50            | U |
| 2,4-dichlorophenol         | 10         | U       | pentachlorophenol     |   | 50            | Ü |
| 2,4-dimethylphenol         | 10         | U       | phenol                |   | 10            | U |
| 2,4-dinitrophenol          | <b>5</b> Ö | U       | 2,4,6-trichlorophenol |   | 10            | U |
| 2-methyl-4,6-dinitrophenol | 50         | U       |                       |   |               |   |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

<sup>-</sup> Indicates an estimated value less than the detection limit.

Client Project ID: LE Carpenter

Job Number: ITAD 51315

### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in  $\mu g$ /liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-063-3 Lab Sample ID: K1756R

| Tentative Identification (1)     | Concentration (2) |
|----------------------------------|-------------------|
| benzothiazole                    | 21                |
| benzothiazole, 2-(methylthio)-   | 9.4               |
| 2-pentanone, 4-hydroxy-4-methyl- | 11 AB             |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- A Suspected aldol condensation product.
- B Analyte was found in the blank as well as the sample.

Job Number: ITAD 51315

### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-063-4 Lab Sample ID:

| Compound                   | Concen | tration  | Compound                    | Concentr    | ation |  |
|----------------------------|--------|----------|-----------------------------|-------------|-------|--|
| acenaphthene               | 500    | U,       | bis(2-chloroisopropyl)ether | 500         | U     |  |
| acenaphthylene             | 500    | U        | bis(2-ethylhexyl)phthalate  | 490         | J     |  |
| anthracene                 | 500    | Ŭ        | 4-bromophenyl phenyl ether  | 500         | Ŭ     |  |
| benzidine                  | 2,500  | U        | 2-chloronaphthalene         | 500         | U     |  |
| benzo(a)anthracene         | 500    | U        | 4-chlorophenyl phenyl ether | 500         | U     |  |
| benzo(b)fluoranthene       | 500    | U        | chrysene                    | 500         | U     |  |
| benzo(k)fluoranthene       | 500    | U        | dibenz(a,h)anthracene       | 500         | U     |  |
| benzo(a)pyrene             | 500    | <b>U</b> | di-n-butylphthalate         | 500         | U     |  |
| benzo(g,h,i)perylene       | 500    | บ        | 1,2-dichlorobenzene         | 500         | U     |  |
| butylbenzylphthalate       | 190    | J        | 1,3-dichlorobenzene         | <b>5</b> 00 | U     |  |
| bis(2-chloroethoxy)methane | 500    | U        | 1,4-dichlorobenzene         | 500         | U.    |  |
| bis(2-chloroethyl)ether    | 500    | U        |                             |             |       |  |

<sup>-</sup> Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction:

05/08/92

Date of Analysis:

05/13/92

Indicates an estimated value less than the detection limit.

Job Number: ITAD 51315

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-063-4 Lab Sample ID: K1757

| Compound                  | Concentra | <u>ation</u> | Compound                   | Concentra | Concentration |  |
|---------------------------|-----------|--------------|----------------------------|-----------|---------------|--|
| 3,3'dichlorobenzidine     | 1,000     | U            | hexachloroethane           | 500       | U             |  |
| diethylphthalate          | 500       | U            | indeno(1,2,3-cd)pyrene     | 500       | Ū             |  |
| dimethyl phthalate        | 500       | Ü            | isophorone                 | 500       | U             |  |
| 2,4-dinitrotoluene        | 500       | U            | naphthalene                | 500       | U             |  |
| 2,6-dinitrotoluene        | 500       | U.           | nitrobenzene               | 500       | U             |  |
| di-n-octyl phthalate      | 500       | Ü            | n-nitroso-di-n-propylamine | 500       | Ü             |  |
| 1,2-diphenylhydrazine(1)  | 500       | U            | n-nitrosodimethylamine     | 500       | U             |  |
| fluoranthene              | 500       | Ŭ            | n-nitrosodiphenylamine(2)  | 500       | U             |  |
| fluorene                  | 500       | U            | phenanthrene               | 500       | U             |  |
| hexachlorobenzene         | 500       | U            | pyrene                     | 500       | U             |  |
| hexachlorobutadiene       | 500       | U            | 1,2,4-trichlorobenzene     | 500       | U             |  |
| hexachlorocyclopentadiene | 500       | U ·          |                            |           |               |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

<sup>(1) -</sup> Screened for as azobenzene

<sup>(2) -</sup> Detected as diphenylamine

Job Number: ITAD 51315

### ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in  $\mu g$ /liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-063-4 Lab Sample ID: K1757

| Compound                   | <u>Concen</u> | tration    | Compound              | Concentration |   |  |
|----------------------------|---------------|------------|-----------------------|---------------|---|--|
| 4-chloro-3-methylphenol    | 500           | U          | 2-nitrophenol         | 500           | U |  |
| 2-chlorophenol             | 500           | U          | 4-nitrophenol         | 2,500         | U |  |
| 2,4-dichlorophenol         | 500           | U          | pentachlorophenol     | 2,500         | Ü |  |
| 2,4-dimethylphenol         | <i>5</i> 00   | U          | phenol                | 500           | U |  |
| 2,4-dinitrophenol          | 2,500         | <b>U</b> , | 2,4,6-trichlorophenol | 500           | U |  |
| 2-methyl-4,6-dinitrophenol | 2,500         | U 🔚        |                       |               |   |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 05/6

05/08/92

Date of Analysis:

05/13/92

J - Indicates an estimated value less than the detection limit.

Job Number: ITAD 51315

### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-063-4 Lab Sample ID: K1757

| Tentative Identification (1)                  | Concentration (2) |
|---|-------------------|
| unknown (organic acid)                        | 9,300             |
| unknown                                       | 2,200             |
| 9-octadecenoic acid (Z)-                      | 1,300             |
| tetradecanoic acid                            | 1,100             |
| hexadecanoic acid                             | 850               |
| 2-pentanone, 4-hydroxy-4-methyl-              | 1,300 AB          |
| dodecanamide, n,n-bis(2-hydroxyethyl)-        | 500               |
| 9-octadecenoic acid (Z)-,2-hydroxyethyl ester | 480               |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- A Suspected aldol condensation product.
- B Analyte was found in the blank as well as the sample.

Job Number: ITAD 51315

#### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 1

Lab Sample ID: H0510R

| Compound                   | Concentration | Compound                    | Concentration |
|----------------------------|---------------|-----------------------------|---------------|
| acenaphthene               | 10 U          | bis(2-chloroisopropyl)ether | 10 U          |
| acenaphthylene             | 10 U          | bis(2-ethylhexyl)phthalate  | 10 U          |
| anthracene                 | 10 U          | 4-bromophenyl phenyl ether  | 10 U          |
| benzidine                  | 50 U          | 2-chloronaphthalene         | 10 U          |
| benzo(a)anthracene         | 10 U          | 4-chlorophenyl phenyl ether | 10 U          |
| benzo(b)fluoranthene       | 10 U          | chrysene                    | 10 U          |
| benzo(k)fluoranthene       | 10 U          | dibenz(a,h)anthracene       | 10 U          |
| benzo(a)pyrene             | 10 U          | di-n-butylphthalate         | 10 U          |
| benzo(g,h,i)perylene       | 10 U          | 1,2-dichlorobenzene         | 10 U          |
| butylbenzylphthalate       | 10 U          | 1,3-dichlorobenzene         | 10 U          |
| bis(2-chloroethoxy)methane | 10 U          | 1,4-dichlorobenzene         | 10 U          |
| bis(2-chloroethyl)ether    | 10 Ú          |                             |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 05/08/92 Date of Analysis: 05/13/92

This method blank applies to the following samples: 1045-063-2 and 1045-063-3.

J - Indicates an estimated value less than the detection limit.

Job Number: ITAD 51315

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 1 Lab Sample ID: H0510R

| Compound                  | Concentration        | Compound                   | Concentration |  |  |
|---------------------------|----------------------|----------------------------|---------------|--|--|
| 3,3'dichlorobenzidine     | 20 U                 | hexachloroethane           | 10 U          |  |  |
| diethylphthalate          | <b>10</b> , <b>U</b> | indeno(1,2,3-cd)pyrene     | 10 · U        |  |  |
| dimethyl phthalate        | 10 <b>Ü</b>          | isophorone                 | 10 U          |  |  |
| 2,4-dinitrotoluene        | 10 U                 | naphthalene                | 10 U          |  |  |
| 2,6-dinitrotoluene        | 10 U                 | nitrobenzene               | 10 U          |  |  |
| di-n-octyl phthalate      | 10 U                 | n-nitroso-di-n-propylamine | 10 U          |  |  |
| 1,2-diphenylhydrazine(1)  | 10 U                 | n-nitrosodimethylamine     | 10 U          |  |  |
| fluoranthene              | 10 U                 | n-nitrosodiphenylamine(2)  | 10 U          |  |  |
| fluorene                  | 10 Ü                 | phenanthrene               | 10 U          |  |  |
| hexachlorobenzene         | 10 Ų                 | pyrene                     | 10 U          |  |  |
| hexachlorobutadiene       | 10 U                 | 1,2,4-trichlorobenzene     | 10 U          |  |  |
| hexachlorocyclopentadiene | 10 U                 |                            |               |  |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

<sup>(1) -</sup> Screened for as azobenzene

<sup>(2) -</sup> Detected as diphenylamine

Job Number: ITAD 51315

### ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 1

Lab Sample ID: H0510R

| Compound                   | Concentration |            | Concentration Compound |    | Concentration |  |  |
|----------------------------|---------------|------------|------------------------|----|---------------|--|--|
| 4-chloro-3-methylphenol    | 10            | Ü          | 2-nitrophenol          | 10 | U             |  |  |
| 2-chlorophenol             | 10            | <b>U</b> : | 4-nitrophenol          | 50 | U             |  |  |
| 2,4-dichlorophenol         | 10            | U          | pentachlorophenol      | 50 | Ù             |  |  |
| 2,4-dimethylphenol         | 10            | U          | phenol                 | 10 | U             |  |  |
| 2,4-dinitrophenol          | 50            | U          | 2,4,6-trichlorophenol  | 10 | U             |  |  |
| 2-methyl-4,6-dinitrophenol | <b>5</b> Ö    | U          |                        |    |               |  |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 05

05/08/92

Date of Analysis:

05/13/92

J - Indicates an estimated value less than the detection limit.

Job Number: ITAD 51315

# ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 1

Lab Sample ID: H0510R

Tentative Identification (1)

Concentration (2)

2-pentanone, 4-hydroxy-4-methyl-

14 A

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- A Suspected aldol condensation product.

Job Number: ITAD 51315

## BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in  $\mu g$ /liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 2

Lab Sample ID: H0516

| Compound                   | Concentration | Compound                    | Concentration |
|----------------------------|---------------|-----------------------------|---------------|
| acenaphthene               | 10 U          | bis(2-chloroisopropyl)ether | 10 U          |
| acenaphthylene             | 10 U          | bis(2-ethylhexyl)phthalate  | 10 U          |
| anthracene                 | 10 U          | 4-bromophenyl phenyl ether  | 10 U          |
| benzidine                  | 50 U          | 2-chloronaphthalene         | 10 U          |
| benzo(a)anthracene         | 10 U          | 4-chlorophenyl phenyl ether | 10 U          |
| benzo(b)fluoranthene       | 10 U          | chrysene                    | 10 Ú          |
| benzo(k)fluoranthene       | 10 U          | dibenz(a,h)anthracene       | 10 U          |
| benzo(a)pyrene             | 10 U          | di-n-butylphthalate         | 10 U          |
| benzo(g,h,i)perylene       | 10 U          | 1,2-dichlorobenzene         | 10 U          |
| butylbenzylphthalate       | 10 U          | 1,3-dichlorobenzene         | 10 U          |
| bis(2-chloroethoxy)methane | 10 U          | 1,4-dichlorobenzene         | 10 Ü          |
| bis(2-chloroethyl)ether    | 10 U          |                             |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 05/08/92 Date of Analysis: 05/13/92

This method blank applies to the following samples: 1045-063-1, 1045-063-1 DL and 1045-063-4.

DL = Dilution

J - Indicates an estimated value less than the detection limit.

Job Number: ITAD 51315

## BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 2

Lab Sample ID: H0516

| Compound                  | Concentration |   | Compound                   | Concentration |            |  |
|---------------------------|---------------|---|----------------------------|---------------|------------|--|
| 3,3'dichlorobenzidine     | 20            | Ŭ | hexachloroethane           | 10            | U          |  |
| diethylphthalate          | 10            | U | indeno(1,2,3-cd)pyrene     | 10            | ·U         |  |
| dimethyl phthalate        | 10            | U | isophorone                 | 10            | U          |  |
| 2,4-dinitrotoluene        | 10            | U | naphthalene                | 10            | U          |  |
| 2,6-dinitrotoluene        | 10            | U | nitrobenzene               | 10            | U          |  |
| di-n-octyl phthalate      | 10            | U | n-nitroso-di-n-propylamine | 10            | U          |  |
| 1,2-diphenylhydrazine(1)  | 10            | U | n-nitrosodimethylamine     | 10            | Ų          |  |
| fluoranthene              | 10            | U | n-nitrosodiphenylamine(2)  | 10            | U          |  |
| fluorene                  | 10            | U | phenanthrene               | 10            | U          |  |
| hexachlorobenzene         | 10            | U | pyrene                     | 10            | U          |  |
| hexachlorobutadiene       | 10            | U | 1,2,4-trichlorobenzene     | 10            | <b>ט</b> , |  |
| hexachlorocyclopentadiene | 10            | U |                            |               |            |  |
|                           |               |   |                            |               |            |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 05/08/92 Date of Analysis: 05/13/92

Indicates an estimated value less than the detection limit.

<sup>(1) -</sup> Screened for as azobenzene

<sup>(2) -</sup> Detected as diphenylamine

Job Number: ITAD 51315

## ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in  $\mu g$ /liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 2

Lab Sample ID: H0516

| Compound                   | Concentration | Compound              | <b>Concentration</b> |
|----------------------------|---------------|-----------------------|----------------------|
| 4-chloro-3-methylphenol    | 10 U          | 2-nitrophenol         | 10 U                 |
| 2-chlorophenol             | 10 U          | 4-nitrophenol         | 50 U                 |
| 2,4-dichlorophenol         | 10 U          | pentachlorophenol     | . <b>50 Ú</b>        |
| 2,4-dimethylphenol         | 10 U          | phenol                | 10 Ų                 |
| 2,4-dinitrophenol          | 50 U          | 2,4,6-trichlorophenol | 10 U                 |
| 2-methyl-4,6-dinitrophenol | 50 U          |                       | •                    |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction:

05/08/92

Date of Analysis:

05/13/92

J - Indicates an estimated value less than the detection limit.

Client Project ID: LE Carpenter

Job Number: ITAD 51315

## ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 2

Lab Sample ID: H0516

Tentative Identification (1)

Concentration (2)

2-pentanone, 4-hydroxy-4-methyl-

28 A

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- A Suspected aldol condensation product.

Job Number: ITAD 51315

## WATER SURROGATE PERCENT RECOVERY SUMMARY

## **SEMI-VOLATILE**

| Client Sample ID | Nitro-<br>Benzene-D5<br>(35-114%)* | 2-Fluoro-<br>Biphenyl<br>(43-116%)* | Terphenyl-<br>D14<br>(33-141%)* | Phenol-D5<br>(10-94%)* | 2-Fluoro-<br>Phenol<br>(21-100%)* | 2,4,6-<br>Tribromo-<br>Phenol<br>(10-123%)* |
|------------------|------------------------------------|-------------------------------------|---------------------------------|------------------------|-----------------------------------|---|
| 1045-063-1       | 76                                 | 86                                  | 160 **                          | 40                     | 57                                | 49  |
| 1045-063-1 DL    | 75                                 | 79                                  | 93                              | 33                     | 49                                | 69  |
| 1045-063-2       | 83                                 | 104                                 | 80                              | 31                     | 47                                | 67  |
| 1045-063-3       | 77                                 | 102                                 | 61                              | 32                     | 46                                | 72  |
| 1045-063-4       | 85                                 | 93                                  | .78                             | 38                     | .53                               | 54  |
| Method Blank 1   | 84                                 | 85                                  | 75                              | 35                     | 52                                | 63  |
| Method Blank 2   | 89                                 | 94                                  | 78                              | 36                     | 53                                | 67  |

DL - Dilution

<sup>\*</sup>Values in parenthesis represent contract required QC limits.

<sup>\*\*</sup>Values are outside of contract required QC limits.

Job Number: ITAD 51315

## CHEMICAL OXYGEN DEMAND

Results in mg/liter (ppm)

Sample Matrix: Water

| Client Sample ID | Lab Sample ID | Result |
|------------------|---------------|--------|
| Method Blank     |               | *      |
| 1045-063-1       | K1750         | 8,700  |
| 1045-063-2       | K1751         | 35     |
| 1045-063-3       | K1752         | 13     |
| 1045-063-4       | K1753         | 9,200  |

Date of Analysis: 05/06/92

<sup>-</sup> A method blank is not applicable for this analysis.

Job Number: ITPK 51491

## **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: 1045-066-01 Lab Sample ID: K3212

| Compound                 | Concentration | on Compound               | Concentra | ation |
|--------------------------|---------------|---------------------------|-----------|-------|
| acrolein                 | <b>50</b> U   | 1,1-dichloroethene        | 25        | U     |
| acrylonitrile            | 50 U          | cis-1,2-dichloroethene    | 25        | U     |
| benzene                  | 16 J          | trans-1,2-dichloroethene  | 25        | U     |
| bromodichloromethane     | 25 U          | 1,2-dichloropropane       | 25        | U     |
| bromoform                | 25 U          | cis-1,3-dichloropropene   | 25        | U     |
| bromomethane             | 50 U          | trans-1,3-dichloropropene | 25        | U     |
| carbon tetrachloride     | 25 U          | ethyl benzene             | 260       |       |
| chlorobenzene            | <b>25</b> U   | methylene chloride        | 1,200     | BD    |
| chloroethane             | 50 U          | 1,1,2,2-tetrachloroethane | 25        | U     |
| 2-chloroethylvinyl ether | 50 U          | tetrachloroethene         | 25        | U     |
| chloroform               | 7 Ј           | toluene                   | 26        |       |
| chloromethane            | 50 U          | 1,1,1-trichloroethane     | 25        | U     |
| dibromochloromethane     | 25 U          | 1,1,2-trichloroethane     | 25        | U     |
| 1,1-dichloroethane       | 25 U          | trichloroethene           | 25        | Ų     |
| 1,2-dichloroethane       | 25 U          | trichlorofluoromethane    | 25        | Ū     |
|                          |               | vinyl chloride            | 50        | U     |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/05/92

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

D - Compound analyzed at a secondary dilution factor.

Job Number: ITPK 51491

## ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-066-01 Lab Sample ID: K3212

| Tentative Identification (1) | Concentration (2) |
|------------------------------|-------------------|
| benzene, dimethyl-           | 510 Y             |
| nonane                       | 22                |
| benzene, 1,2-dimethyl        | 31                |
| unknown (SAT'D HC)           | 29                |
| nonane, 3-methyl-            | 26                |
| benzene, ethylmethyl-        | 20 Y              |
| decane                       | . 58              |
| benzene, trimethyl-          | 22 Y              |
| decane, 4-methyl-            | 20                |
| undecane                     | 240               |

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Y - Indistinguishable isomer in tentatively identified compounds.

Job Number: ITPK 51491

## **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: 1045-066-02 Lab Sample ID: K3213R

| Compound                 | Concentration | Compound                  | Concent | ration |
|--------------------------|---------------|---------------------------|---------|--------|
| acrolein                 | 33 U          | 1,1-dichloroethene        | 17      | U      |
| acrylonitrile            | 33 U          | cis-1,2-dichloroethene    | 17      | U      |
| benzene                  | 6 J           | trans-1,2-dichloroethene  | 17      | U      |
| bromodichloromethane     | 17 U          | 1,2-dichloropropane       | 17      | U      |
| bromoform                | 17 U          | cis-1,3-dichloropropene   | 17      | U      |
| bromomethane             | 33 U          | trans-1,3-dichloropropene | · 17    | U      |
| carbon tetrachloride     | 17 U          | ethyl benzene             | 17      | U      |
| chlorobenzene            | 17 U          | methylene chloride        | 390     | В      |
| chloroethane             | 33 U          | 1,1,2,2-tetrachloroethane | 17      | U      |
| 2-chloroethylvinyl ether | 33 Ū          | tetrachloroethene         | 17      | U      |
| chloroform               | 17 U          | toluene                   | 16      | j .    |
| chloromethane            | 33 U          | 1,1,1-trichloroethane     | 17      | U      |
| dibromochloromethane     | 17 U          | 1,1,2-trichloroethane     | 17      | U      |
| 1,1-dichloroethane       | 17 U          | trichloroethene           | 17      | U      |
| 1,2-dichloroethane       | 17 U          | trichlorofluoromethane    | 17      | U      |
|                          | • .           | vinyl chloride            | 33      | U      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/05/92

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

Client Project ID: LE Carpenter

Job Number: ITPK 51491

## ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-066-02 Lab Sample ID: K3213R

| Tentative Identification (1) | Concentration (2) |
|------------------------------|-------------------|
| benzene, dimethyl-           | 54 Y              |
| benzene, 1,2-dimethyl-       | 11                |

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Y - Indistinguishable isomer in tentatively identified compounds.

Job Number: ITPK 51491

## **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: 1045-066-03 Lab Sample ID: K3214R

| Compound                 | Concentration | Compound                  | Concent | ration |
|--------------------------|---------------|---------------------------|---------|--------|
| acrolein                 | 10 ° U        | 1,1-dichloroethene        | 5       | U      |
| acrylonitrile            | 10 U          | cis-1,2-dichloroethene    | 5       | U      |
| benzene                  | 2 J           | trans-1,2-dichloroethene  | 5       | U      |
| bromodichloromethane     | 1 J           | 1,2-dichloropropane       | 5       | Ų      |
| bromoform                | 5 U           | cis-1,3-dichloropropene   | 5       | U      |
| bromomethane             | 10 U          | trans-1,3-dichloropropene | . 5     | U      |
| carbon tetrachloride     | 5 U           | ethyl benzene             | 3       | J      |
| chlorobenzene            | 5 U           | methylene chloride        | 5,800   | BD     |
| chloroethane             | 10 U          | 1,1,2,2-tetrachloroethane | 5       | U      |
| 2-chloroethylvinyl ether | 10 U          | tetrachloroethene         | 5       | U      |
| chloroform               | 18            | toluene                   | 3       | J      |
| chloromethane            | 10 U          | 1,1,1-trichloroethane     | 5       | U      |
| dibromochloromethane     | <b>5 U</b>    | 1,1,2-trichloroethane     | 5       | U      |
| 1,1-dichloroethane       | 5 U           | trichloroethene           | 5       | U      |
| 1,2-dichloroethane       | 5 U           | trichlorofluoromethane    | 5       | U      |
|                          |               | vinyl chloride            | 10      | U      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/08/92

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

D - Compound analyzed at a secondary dilution factor.

Client Project ID: LE Carpenter

Job Number: ITPK 51491

## ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-066-03 Lab Sample ID: K3214R

| Tentative Identification (1) | Concentra | tion ( | <u>2)</u> |
|------------------------------|-----------|--------|-----------|
| benzene, dimethyl-           | .28       | Y      |           |
| benzene, 1,2-dimethyl-       | 7.6       |        | ,         |

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Y - Indistinguishable isomer in tentatively identified compounds.

Job Number: ITPK 51491

## VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)
Sample Matrix: Water

Client Sample ID: 1045-066-04 Lab Sample ID: K3215

| Compound                 | Concentrati | <u>ion</u> | Compound                  | Concent | ration |
|--------------------------|-------------|------------|---------------------------|---------|--------|
| acrolein                 | 50 U        | J          | 1,1-dichloroethene        | 25      | U      |
| acrylonitrile            | 50 U        | J          | cis-1,2-dichloroethene    | 25      | U      |
| benzene                  | 15 J        | ,          | trans-1,2-dichloroethene  | 25      | Ü      |
| bromodichloromethane     | 25 U        | J' :       | 1,2-dichloropropane       | 25      | U      |
| bromoform                | 25 U        | J          | cis-1,3-dichloropropene   | 25      | Ù      |
| bromomethane             | 50 U        | J          | trans-1,3-dichloropropene | 25      | U      |
| carbon tetrachloride     | 25 U        | J .        | ethyl benzene             | 25      | U      |
| chlorobenzene            | 25 U        | J          | methylene chloride        | 58      | В      |
| chloroethane             | 50 U        | J          | 1,1,2,2-tetrachloroethane | 25      | U      |
| 2-chloroethylvinyl ether | 50 U        | j          | tetrachioroethene         | 25      | U      |
| chloroform               | 9 J         |            | toluene                   | 21      | j      |
| chloromethane            | 50 Ü        | Ť .        | 1,1,1-trichloroethane     | 25      | U      |
| dibromochloromethane     | 25 U        | J          | 1,1,2-trichloroethane     | 25      | U      |
| 1,1-dichloroethane       | 25 U        | J          | trichloroethene           | 25      | U      |
| 1,2-dichloroethane       | 25 U        | Ţ.         | trichlorofluoromethane    | 25      | Ü      |
|                          |             | :          | vinyl chloride            | 50      | ,<br>T |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/08/92

<sup>-</sup> Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

Client Project ID: LE Carpenter

Job Number: ITPK 51491

## ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-066-04 Lab Sample ID: K3215

| Tentative Identification (1) | Concentration (2) |
|------------------------------|-------------------|
| pentane                      | 100               |
| hexane                       | 110               |
| octane                       | 76                |
| benzene, dimethyl-           | 42 Y              |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- Y Indistinguishable isomer in tentatively identified compounds.

Job Number: ITPK 51491

## CHEMICAL OXYGEN DEMAND ANALYSIS

Results in mg/liter (ppm)

Sample Matrix: Water

| Client Sample ID | Lab Sample ID | <u>Result</u> | Analysis Date |
|------------------|---------------|---------------|---------------|
| Method Blank     | *             | -             |               |
| 1045-066-01      | K3224         | 11,000        | 06/02/92      |
| 1045-066-02      | K3225         | 34            | 06/03/92      |
| 1045-066-03      | K3226         | 7.5           | 06/03/92      |
| 1045-066-04      | K3227         | 11,000        | 06/02/92      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

<sup>\* -</sup> A method blank is not applicable for this analysis.

Job Number: ITPK 51491

## **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in µg/liter (ppb)
Sample Matrix: Water

Client Sample ID: 1045-068-01 Lab Sample ID: K3216

| Compound                 | Concent | ration | Compound                  | Concent | ration |
|--------------------------|---------|--------|---------------------------|---------|--------|
| acrolein                 | 10      | U      | 1,1-dichloroethene        | 5       | U      |
| acrylonitrile            | 10      | U      | cis-1,2-dichloroethene    | 5       | U      |
| benzene                  | 26      |        | trans-1,2-dichloroethene  | 5       | U      |
| bromodichloromethane     | 3       | J      | 1,2-dichloropropane       | 5       | U      |
| bromoform                | 5       | Ū      | cis-1,3-dichloropropene   | 5       | U      |
| bromomethane             | 10      | U      | trans-1,3-dichloropropene | 5       | U      |
| carbon tetrachloride     | 5       | Ü      | ethyl benzene             | 7       |        |
| chlorobenzene            | 5       | U      | methylene chloride        | 1,300   | BD     |
| chloroethane             | 10      | Ū      | 1,1,2,2-tetrachloroethane | 5       | U      |
| 2-chloroethylvinyl ether | 10      | U      | tetrachloroethene         | 5       | U      |
| chloroform               | 32      |        | toluene                   | 29      | •      |
| chloromethane            | 10      | Ū      | 1,1,1-trichloroethane     | 5       | U      |
| dibromochloromethane     | 5       | U      | 1,1,2-trichloroethane     | 5       | U      |
| 1,1-dichloroethane       | 5       | U      | trichloroethene           | . 5     | U      |
| 1,2-dichloroethane       | 5       | U      | trichlorofluoromethane    | 5       | U      |
|                          |         | •      | vinyl chloride            | 10      | U      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/09/92

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

D - Compound analyzed at a secondary dilution factor.

Job Number: ITPK 51491

## ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-068-01 Lab Sample ID: K3216

| Tentative Identification (1) | Concentra | tion (2)  |
|------------------------------|-----------|-----------|
| carbon disulfide             |           |           |
|                              | 20        |           |
| 2-propenoic acid, 2-methyl-, | 5.3       |           |
| benzene, dimethyl-           | 2,700     | Y         |
| benzene, 1,2-dimethyl-       | 560       |           |
| benzene, ethylmethyl-        | 40        | <b>'Y</b> |
| benzene, trimethyl-          | 39        | Y         |
| benzene, ethylmethyl-        | 7.2       | Y         |
| benzene, trimethyl-          | 58        | Y         |
| benzene, trimethyl-          | 24        | Y         |
| benzene, tetramethyl-        | 8.1       | Ÿ         |

<sup>(1)</sup> Identification is based on computer search of the NIST Library.

<sup>(2)</sup> Concentration is based on a response factor of 1.00 relative to the internal standard.

Y - Indistinguishable isomer in tentatively identified compounds.

Job Number: ITPK 51491

## **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: 1045-068-02 Lab Sample ID: K3217

| Compound                 | Concent    | ration | Compound                  | Concent   | ration |
|--------------------------|------------|--------|---------------------------|-----------|--------|
| acrolein                 | 10         | U      | 1,1-dichloroethene        | <b>.5</b> | U      |
| acrylonitrile            | 10         | ប      | cis-1,2-dichloroethene    | -5        | U      |
| benzene                  | .3         | J      | trans-1,2-dichloroethene  | 5         | Ü      |
| bromodichloromethane     | 2          | J      | 1,2-dichloropropane       | 5         | U      |
| bromoform                | · <b>5</b> | U      | cis-1,3-dichloropropene   | 5         | U      |
| bromomethane             | 10         | Ū      | trans-1,3-dichloropropene | 5         | U      |
| carbon tetrachloride     | 5          | U      | ethyl benzene             | . 4       | J      |
| chlorobenzene            | 5          | Ü      | methylene chloride        | 410       | BD     |
| chloroethane             | 10         | ប      | 1,1,2,2-tetrachloroethane | 5         | U      |
| 2-chloroethylvinyl ether | 10         | U      | tetrachloroethene         | 5         | U      |
| chloroform               | 20         | č      | toluene                   | 5         |        |
| chloromethane            | 10         | U      | 1,1,1-trichloroethane     | 5         | U      |
| dibromochloromethane     | 5          | U      | 1,1,2-trichloroethane     | 5         | U      |
| 1,1-dichloroethane       | 5          | ប      | trichloroethene           | 5         | U      |
| 1,2-dichloroethane       | 5          | U      | trichlorofluoromethane    | 5         | U      |
|                          |            | 1      | vinyl chloride            | 10        | U      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/09/92

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

D - Compound analyzed at a secondary dilution factor.

Client Project ID: LE Carpenter

Job Number: ITPK 51491

## ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-068-02 Lab Sample ID: K3217

Tentative Identification (1)

benzene, dimethyl
benzene, 1,2-dimethyl
13

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- Y Indistinguishable isomer in tentatively identified compounds.

Job Number: ITPK 51491

## **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu$ g/liter (ppb) Sample Matrix: Water

Client Sample ID: 1045-068-03 Lab Sample ID: K3218

| Compound                 | Concentration     | Compound                  | Concent | ration |
|--------------------------|-------------------|---------------------------|---------|--------|
| acrolein                 | 10 † <b>U</b>     | 1,1-dichloroethene        | 5       | U      |
| acrylonitrile            | 10 Ú              | cis-1,2-dichloroethene    | 5       | U      |
| benzene                  | 3. J              | trans-1,2-dichloroethene  | 5       | Ü      |
| bromodichloromethane     | 3 Ј               | 1,2-dichloropropane       | 5       | U      |
| bromoform                | 5 🖁 บ             | cis-1,3-dichloropropene   | 5       | U      |
| bromomethane             | 10 <sup>°</sup> U | trans-1,3-dichloropropene | 5       | U      |
| carbon tetrachloride     | 5 U               | ethyl benzene             | 2       | J      |
| chlorobenzene            | 5 U               | methylene chloride        | 5,600   | BD     |
| chloroethane             | 10 U              | 1,1,2,2-tetrachloroethane | 5       | U      |
| 2-chloroethylvinyl ether | 10 U              | tetrachloroethene         | 5       | U      |
| chloroform               | 26                | toluene                   | 5       |        |
| chloromethane            | 10 Ú              | 1,1,1-trichloroethane     | 5       | U      |
| dibromochloromethane     | 5 U               | 1,1,2-trichloroethane     | 5       | U      |
| 1,1-dichloroethane       | 5 U               | trichloroethene           | 5       | U      |
| 1,2-dichloroethane       | 5 U               | trichlorofluoromethane    | 5       | U      |
|                          |                   | vinyl chloride            | 10      | U.     |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/09/92

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

D - Compound analyzed at a secondary dilution factor.

Client Project ID: LE Carpenter

Job Number: ITPK 51491

## ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-068-03 Lab Sample ID: K3218

| Tentative Identification (1) | Concentration (2) |
|------------------------------|-------------------|
| benzene, dimethyl-           | 28 Y              |
| benzene, 1,2-dimethyl-       | <b>5.5</b> '      |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- Y Indistinguishable isomer in tentatively identified compounds.

Job Number: ITPK 51491

## **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: 1045-068-04 Lab Sample ID: K3219

| Compound                 | Concentra    | <u>tion</u> | Compound                  | Concent | ration |
|--------------------------|--------------|-------------|---------------------------|---------|--------|
| acrolein                 | 10           | U           | 1,1-dichloroethene        | 5       | U      |
| acrylonitrile            | 10 l         | U           | cis-1,2-dichloroethene    | 5       | U      |
| benzene                  | <i>5</i> (   | U           | trans-1,2-dichloroethene  | 5       | U      |
| bromodichloromethane     | 3 J          | J           | 1,2-dichloropropane       | 5       | U      |
| bromoform                | <b>5</b> , t | <b>U</b> ,  | cis-1,3-dichloropropene   | 5       | U      |
| bromomethane             | 10 U         | ָ<br>ט      | trans-1,3-dichloropropene | 5       | U      |
| carbon tetrachloride     | 5 T          | ע           | ethyl benzene             | 1       | J      |
| chlorobenzene            | 5 T          | ט           | methylene chloride        | 53      | В      |
| chloroethane             | 10 T         | ט           | 1,1,2,2-tetrachloroethane | 5       | Ü      |
| 2-chloroethylvinyl ether | 10 T         | U           | tetrachloroethene         | 5       | U      |
| chloroform               | 28           |             | toluene                   | 2       | J      |
| chloromethane            | 10 T         | U           | 1,1,1-trichloroethane     | 1       | J      |
| dibromochloromethane     | 5 T          | U           | 1,1,2-trichloroethane     | . 5     | U      |
| 1,1-dichloroethane       | 5 τ          | ע           | trichloroethene           | ·<br>5  | U      |
| 1,2-dichloroethane       | 5 T          | ט           | trichlorofluoromethane    | 5       | U      |
|                          |              |             | vinyl chloride            | 10      | U      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/09/92

Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

IT Corporation June 23, 1992 IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

Client Project ID: LE Carpenter

Job Number: ITPK 51491

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-068-04 Lab Sample ID: K3219

Tentative Identification (1)

Concentration (2)

None Detected

Remarks:

(1) Identification is based on computer search of the NIST Library.

(2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITPK 51491

## **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: Method Blank 1 Lab Sample ID: WB06052

| Compound                 | <u>Concentration</u> | Compound                  | Concentration |
|--------------------------|----------------------|---------------------------|---------------|
| acrolein                 | 10 U                 | 1,1-dichloroethene        | 5 U           |
| acrylonitrile            | 10 U                 | cis-1,2-dichloroethene    | 5 U           |
| benzene                  | 5 °U                 | trans-1,2-dichloroethene  | 5 U           |
| bromodichloromethane     | 5 U                  | 1,2-dichloropropane       | 5 U           |
| bromoform                | 5 U                  | cis-1,3-dichloropropene   | 5 U           |
| bromomethane             | 10 Ü                 | trans-1,3-dichloropropene | . 5 U         |
| carbon tetrachloride     | .5 U                 | ethyl benzene             | 5 Ü           |
| chlorobenzene            | 5 Ü                  | methylene chloride        | 2 Ј           |
| chloroethane             | 10 U                 | 1,1,2,2-tetrachloroethane | 5 U           |
| 2-chloroethylvinyl ether | 10 U                 | tetrachloroethene         | 5 U           |
| chloroform               | 5 U                  | toluene                   | 5 U           |
| chloromethane            | 10 U                 | 1,1,1-trichloroethane     | 5 U           |
| dibromochloromethane     | 5 U                  | 1,1,2-trichloroethane     | 5 U           |
| 1,1-dichloroethane       | 5 U                  | trichloroethene           | 5 U           |
| 1,2-dichloroethane       | <b>5</b> U           | trichlorofluoromethane    | 5 U           |
|                          | . "                  | vinyl chloride            | 10 U          |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/05/92

This method blank applies to the following samples: 1045-066-01, 1045-066-01 DL, 1045-066-02.

DL - Dilution.

Indicates an estimated value less than the detection limit.

Client Project ID: LE Carpenter

Job Number: ITPK 51491

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 1 Lab Sample ID: WB06052

Tentative Identification (1)

Concentration (2)

None Detected

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITPK 51491

## **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: Method Blank 2 Lab Sample ID: WB06084

| Compound                 | Concent | ration       | Compound                  | Concent | ration |
|--------------------------|---------|--------------|---------------------------|---------|--------|
| acrolein                 | 10      | U            | 1,1-dichloroethene        | 5       | U      |
| acrylonitrile            | 10      | Ü            | cis-1,2-dichloroethene    | 5       | U      |
| benzene                  | 5       | U            | trans-1,2-dichloroethene  | 5       | U      |
| bromodichloromethane     | 5       | U            | 1,2-dichloropropane       | 5       | U      |
| bromoform                | 5       | U            | cis-1,3-dichloropropene   | 5       | U      |
| bromomethane             | 10      | <b>u</b> , ' | trans-1,3-dichloropropene | 5.      | U      |
| carbon tetrachloride     | 5       | U            | ethyl benzene             | 5       | U      |
| chlorobenzene            | 5       | U            | methylene chloride        | 2       | J      |
| chloroethane             | 10      | U .          | 1,1,2,2-tetrachloroethane | 5       | U      |
| 2-chloroethylvinyl ether | 10      | U            | tetrachloroethene         | 5       | U      |
| chloroform               | 5       | U            | toluene                   | . 5     | U .    |
| chloromethane            | 10      | U            | 1,1,1-trichloroethane     | 5       | U      |
| dibromochloromethane     | 5       | <b>U</b>     | 1,1,2-trichloroethane     | 5       | U      |
| 1,1-dichloroethane       | 5       | U            | trichloroethene           | . 5     | U      |
| 1,2-dichloroethane       | 5       | U            | trichlorofluoromethane    | 5       | U      |
|                          |         |              | vinyl chloride            | 10      | U      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/08/92

This method blank applies to the following samples: 1045-066-03, 1045-066-03 DL, 1045-066-04 and 1045-068-01.

DL - Dilution.

J - Indicates an estimated value less than the detection limit.

Client Project ID: LE Carpenter

Job Number: ITPK 51491

## ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 2 Lab Sample ID: WB06084

Tentative Identification (1)

Concentration (2)

None Detected

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITPK 51491

## **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: Method Blank 3 Lab Sample ID: WB0609

| Compound                 | Concent | ration     | Compound                  | Concent | ration |
|--------------------------|---------|------------|---------------------------|---------|--------|
| acrolein                 | 10      | U          | 1,1-dichloroethene        | 5       | U      |
| acrylonitrile            | 10      | Ū          | cis-1,2-dichloroethene    | 5       | U      |
| benzene                  | 5       | U          | trans-1,2-dichloroethene  | 5       | U      |
| bromodichloromethane     | 5       | U          | 1,2-dichloropropane       | 5       | U      |
| bromoform                | 5       | U          | cis-1,3-dichloropropene   | 5       | Ù      |
| bromomethane             | 10      | U          | trans-1,3-dichloropropene | 5       | U      |
| carbon tetrachloride     | 5       | U          | ethyl benzene             | 5       | U      |
| chlorobenzene            | 5       | U          | methylene chloride        | 2       | . 1    |
| chloroethane             | 10      | Ū          | 1,1,2,2-tetrachloroethane | 5       | U      |
| 2-chloroethylvinyl ether | 10      | U          | tetrachloroethene         | 5       | U      |
| chloroform               | 5       | <b>, U</b> | toluene                   | 5       | Ŭ      |
| chloromethane            | 10      | Ū          | 1,1,1-trichloroethane     | 5       | U      |
| dibromochloromethane     | 5       | Ü          | 1,1,2-trichloroethane     | 5       | U      |
| 1,1-dichloroethane       | 5       | U          | trichloroethene           | 5       | Ù      |
| 1,2-dichloroethane       | 5       | U          | trichlorofluoromethane    | 5       | U      |
|                          |         | :          | vinyl chloride            | 10      | U      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/09/92

This method blank applies to the following samples: 1045-068-01 DL, 1045-068-02, 1045-068-02 DL, 1045-068-03, 1045-068-03 DL, 1045-068-04.

DL - Dilution.

J - Indicates an estimated value less than the detection limit.

Client Project ID: LE Carpenter

Job Number: ITPK 51491

## ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in  $\mu g$ /liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 3

Lab Sample ID: WB0609

Tentative Identification (1)

None Detected

Concentration (2)

Remarks:

(1) Identification is based on computer search of the NIST Library.

(2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITPK 51491

## WATER SURROGATE PERCENT RECOVERY SUMMARY

|                  | 1                        | VOLATILE          |                                     |
|------------------|--------------------------|-------------------|-------------------------------------|
| Client Sample ID | Toluene-D8<br>(88-110%)* | BFB<br>(86-115%)* | 1,2 Dichloroethane-D4<br>(76-114%)* |
| 1045-066-01      | 100                      | 107               | 101                                 |
| 1045-066-01 DL   | 99                       | 96                | 96                                  |
| 1045-066-02      | 96                       | 93                | 93                                  |
| 1045-066-03      | 103                      | 95                | 88                                  |
| 1045-066-03 DL   | 102                      | 98                | 95                                  |
| 1045-066-04      | 106                      | 89                | 96                                  |
| 1045-068-01      | 101                      | 121 ***           | 90                                  |
| 1045-068-01 DL   | 104                      | 109               | 100                                 |
| 1045-068-02      | 98                       | 98                | 92                                  |
| 1045-068-02 DL   | 100                      | 98                | 94                                  |
| 1045-068-03      | 98                       | 99                | 92                                  |
| 1045-068-03 DL   | 104                      | 100               | 97                                  |
| 1045-068-04      | 96                       | 96                | 91                                  |
| Method Blank 1   | 101                      | 97                | 98                                  |
| Method Blank 2   | 100                      | 96                | 95                                  |
| Method Blank 3   | 108                      | 105               | 101                                 |

<sup>\*</sup>Values in parenthesis represent QC limits.

\*\*Values are outside of QC limits.

DL - Dilution

Job Number: ITPK 51491

## BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-066-01 Lab Sample ID: K3232

| Compound                   | Concentration | Compound                    | Concentration |
|----------------------------|---------------|-----------------------------|---------------|
| acenaphthene               | 1,000 U       | bis(2-chloroisopropyl)ether | 1,000 U       |
| acenaphthylene             | 1,000 U       | bis(2-ethylhexyl)phthalate  | 3,200         |
| anthracene                 | 1,000 U       | 4-bromophenyl phenyl ether  | 1,000 U       |
| benzidine                  | 5,000 U       | 2-chloronaphthalene         | 1,000 Ü       |
| benzo(a)anthracene         | 1,000 U       | 4-chlorophenyl phenyl ether | 1,000 U       |
| benzo(b)fluoranthene       | 1,000 U       | chrysene                    | 1,000 U       |
| benzo(k)fluoranthene       | 1,000 U       | dibenz(a,h)anthracene       | 1,000 U       |
| benzo(a)pyrene             | 1,000 U       | di-n-butylphthalate         | 1,000 U       |
| benzo(g,h,i)perylene       | 1,000 U       | 1,2-dichlorobenzene         | 1,000 Ü       |
| butylbenzylphthalate       | 1,000 U       | 1,3-dichlorobenzene         | 1,000 U       |
| bis(2-chloroethoxy)methane | 1,000 U       | 1,4-dichlorobenzene         | 1,000 U       |
| bis(2-chloroethyl)ether    | 1,000 U       |                             | •             |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 06/02/92 Date of Analysis: 06/09/92

J - Indicates an estimated value less than the detection limit.

Job Number: ITPK 51491

## BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-066-01 Lab Sample ID: K3232

| Compound                  | Concentration |   | Compound                   | Concentration |   |
|---------------------------|---------------|---|----------------------------|---------------|---|
| 3,3'dichlorobenzidine     | 2,000         | U | hexachloroethane           | 1,000         | Ŭ |
| diethylphthalate          | 1,000         | Ü | indeno(1,2,3-cd)pyrene     | 1,000         | U |
| dimethyl phthalate        | 1,000         | Ü | isophorone                 | 1,000         | U |
| 2,4-dinitrotoluene        | 1,000         | U | naphthalene                | 1,000         | Ü |
| 2,6-dinitrotoluene        | 1,000         | U | nitrobenzene               | 1,000         | U |
| di-n-octyl phthalate      | 1,000         | U | n-nitroso-di-n-propylamine | 1,000         | U |
| 1,2-diphenylhydrazine(1)  | 1,000         | U | n-nitrosodimethylamine     | 1,000         | U |
| fluoranthene              | 1,000         | Ü | n-nitrosodiphenylamine(2)  | 1,000         | U |
| fluorene                  | 1,000         | U | phenanthrene               | 1,000         | U |
| hexachlorobenzene         | 1,000         | U | ругеле                     | 1,000         | Ù |
| hexachlorobutadiene       | 1,000         | U | 1,2,4-trichlorobenzene     | 1,000         | U |
| hexachlorocyclopentadiene | 1,000         | U |                            |               |   |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 06/02/92 Date of Analysis: 06/09/92

J - Indicates an estimated value less than the detection limit.

<sup>(1) -</sup> Screened for as azobenzene

<sup>(2) -</sup> Detected as diphenylamine

Job Number: ITPK 51491

## ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-066-01 Lab Sample ID: K3232

| Compound                   | Concentration   | Compound              | Concentration |
|----------------------------|-----------------|-----------------------|---------------|
| 4-chloro-3-methylphenol    | 1, <b>000</b> U | 2-nitrophenol         | 1,000 U       |
| 2-chlorophenol             | 1,000 U         | 4-nitrophenol         | 5,000 U       |
| 2,4-dichlorophenol         | 1,000 U         | pentachlorophenol     | 5,000 U       |
| 2,4-dimethylphenol         | 1,000 Ū         | phenol                | 1,000 U       |
| 2,4-dinitrophenol          | 5,000 U         | 2,4,6-trichlorophenol | 1,000 U       |
| 2-methyl-4,6-dinitrophenol | <b>5,000</b> U  |                       | *             |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 06/02/92 Date of Analysis: 06/09/92

J - Indicates an estimated value less than the detection limit.

Job Number: ITPK 51491

## ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-066-01 Lab Sample ID: K3232

| Tentative Identification (1)    | Concentration (2) |
|---------------------------------|-------------------|
| unknown                         | 34,000            |
| 2-hexadecanol                   | 33,000            |
| ethanol, 2-(dodecyloxy)-        | 32,000            |
| unknown                         | 56,000            |
| 1-dodecanol                     | 27,000            |
| unknown                         | 26,000            |
| 1-tetradecanol                  | 26,000            |
| ethanol, 2-(dodecyloxy)-        | 25,000            |
| unknown                         | 43,000            |
| unknown                         | 41,000            |
| unknown                         | 28,000            |
| 1-hexadecanol                   | 9,400             |
| ethanol, 2-(hexadecyloxy)-      | 7,300             |
| unknown                         | 11,000            |
| unknown                         | 9,000             |
| unknown                         | 7,700             |
| unknown                         | 4,200             |
| unknown                         | 6,300             |
| unknown                         | 4,900             |
| cyclododecane                   | 1,900             |
| unknown                         | 2,100             |
| cyclododecanol                  | 1,600             |
| 1-hexadecanol                   | 1,500             |
| unknown                         | 2,400             |
| decanedioic acid, didecyl ester | 1,000             |
|                                 |                   |

<sup>(1)</sup> Identification is based on computer search of the NIST Library.

<sup>(2)</sup> Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITPK 51491

## BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-066-02 Lab Sample ID: K3233

| Compound                   | Concentration | Compound                    | Concentration |  |
|----------------------------|---------------|-----------------------------|---------------|--|
| acenaphthene               | 10 U          | bis(2-chloroisopropyl)ether | 10 U          |  |
| acenaphthylene             | 10 U          | bis(2-ethylhexyl)phthalate  | 10 Ú          |  |
| anthracene                 | 10 U          | 4-bromophenyl phenyl ether  | 10 U          |  |
| benzidine                  | 50 U          | 2-chloronaphthalene         | 10 U          |  |
| benzo(a)anthracene         | 10 U          | 4-chlorophenyl phenyl ether | 10 U          |  |
| benzo(b)fluoranthene       | 10 U          | chrysene                    | . 10 Ü        |  |
| benzo(k)fluoranthene       | 10 U          | dibenz(a,h)anthracene       | 10 U          |  |
| benzo(a)pyrene             | 10 · U        | di-n-butylphthalate         | 10 U          |  |
| benzo(g,h,i)perylene       | 10 U          | 1,2-dichlorobenzene         | 10 U          |  |
| butylbenzylphthalate       | 10 U          | 1,3-dichlorobenzene         | 10 U          |  |
| bis(2-chloroethoxy)methane | 10 U          | 1,4-dichlorobenzene         | 10 Ü          |  |
| bis(2-chloroethyl)ether    | 10 U          |                             |               |  |

Date of Extraction: 06/02/92 Date of Analysis:

06/08/92

Compound was analyzed for but not detected. The number is the detection limit for the sample.

Indicates an estimated value less than the detection limit.

Job Number: ITPK 51491

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-066-02 Lab Sample ID: K3233

| Compound                  | Concentration | Compound                               | Concentration |  |
|---------------------------|---------------|--|---------------|--|
| 3,3'dichlorobenzidine     | 20 U          | hexachloroethane                       | 10 U          |  |
| diethylphthalate          | 2 J           | indeno(1,2,3-cd)pyrene                 | 10 U          |  |
| dimethyl phthalate        | 10 U          | isophorone                             | 10 U          |  |
| 2,4-dinitrotoluene        | 10 U          | naphthalene                            | 10 U          |  |
| 2,6-dinitrotoluene        | 10 <b>U</b>   | nitrobenzene                           | 10 U          |  |
| di-n-octyl phthalate      | 10 U          | n-nitroso-di-n-propylamine             | 10 U          |  |
| 1,2-diphenylhydrazine(1)  | 10 U          | n-nitrosodimethylamine                 | 10 U          |  |
| fluoranthene              | 10 U          | n-nitrosodiphenylamine(2)              | 10 Ü          |  |
| fluorene                  | 10 U          | phenanthrene                           | 10 U          |  |
| hexachlorobenzene         | 10 Ü          | pyrene                                 | 10 U          |  |
| hexachlorobutadiene       | 10 U          | 1,2,4-trichlorobenzene                 | 10 U          |  |
| hexachlorocyclopentadiene | 10 U          | * ************************************ | ·             |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 06/02/92 Date of Analysis: 06/08/92

J - Indicates an estimated value less than the detection limit.

<sup>(1) -</sup> Screened for as azobenzene

<sup>(2) -</sup> Detected as diphenylamine

Job Number: ITPK 51491

#### ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in  $\mu g$ /liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-066-02 Lab Sample ID: K3233

| Compound                   | Concentration | Compound              | Concentration |
|----------------------------|---------------|-----------------------|---------------|
| 4-chloro-3-methylphenol    | 10 U          | 2-nitrophenol         | 10 U          |
| 2-chlorophenol             | 10 U          | 4-nitrophenol         | 50 U          |
| 2,4-dichlorophenol         | 10 U          | pentachlorophenol     | 50 U          |
| 2,4-dimethylphenol         | 10 U          | phenol                | 10 U          |
| 2,4-dinitrophenol          | 50 U          | 2,4,6-trichlorophenol | 10 U          |
| 2-methyl-4,6-dinitrophenol | <b>50</b> U   |                       |               |

U Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction:

06/02/92

Date of Analysis:

06/08/92

Indicates an estimated value less than the detection limit.

Client Project ID: LE Carpenter

Job Number: ITPK 51491

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-066-02 Lab Sample ID: K3233

| Tentative Identification (1)   | Concentration (2) |
|--------------------------------|-------------------|
| 2-pentanone, 4-hydroxy-4-met   | 22 AB             |
| benzothiazole, 2-(methylthio)- | 16                |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- A Suspected aldol condensation product.
- B Analyte was found in the blank as well as the sample.

Job Number: ITPK 51491

#### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-066-03 Lab Sample ID: K3234

| Compound                   | Concen | tration | Compound                    | Concentr | ation |
|----------------------------|--------|---------|-----------------------------|----------|-------|
| acenaphthene               | 10     | U ···   | bis(2-chloroisopropyl)ether | 10       | U     |
| acenaphthylene             | 10     | U       | bis(2-ethylhexyl)phthalate  | . 10     | U     |
| anthracene                 | 10     | U       | 4-bromophenyl phenyl ether  | 10       | U     |
| benzidine                  | 50     | U       | 2-chloronaphthalene         | 10       | U     |
| benzo(a)anthracene         | 10     | U       | 4-chlorophenyl phenyl ether | 10       | U     |
| benzo(b)fluoranthene       | 10     | U       | chrysene                    | 10       | U     |
| benzo(k)fluoranthene       | 10     | U       | dibenz(a,h)anthracene       | 10       | U     |
| benzo(a)pyrene             | 10     | U,      | di-n-butylphthalate         | 10       | U     |
| benzo(g,h,i)perylene       | 10     | U       | 1,2-dichlorobenzene         | 10       | U     |
| butylbenzylphthalate       | 10     | U       | 1,3-dichlorobenzene         | 10       | U     |
| bis(2-chloroethoxy)methane | 10     | U       | 1,4-dichlorobenzene         | 10       | U     |
| bis(2-chloroethyl)ether    | 10     | U ,     |                             |          | *     |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Job Number: ITPK 51491

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-066-03 Lab Sample ID: K3234

| Compound                  | Concentration | Compound                   | Concentration |
|---------------------------|---------------|----------------------------|---------------|
| 3,3'dichlorobenzidine     | 20 U          | hexachloroethane           | 10 U          |
| diethylphthalate          | 10 U          | indeno(1,2,3-cd)pyrene     | 10 U          |
| dimethyl phthalate        | 10 U          | isophorone                 | 10 U          |
| 2,4-dinitrotoluene        | 10 Ü          | naphthalene                | 10 U          |
| 2,6-dinitrotoluene        | 10 Ü          | nitrobenzene               | 10 U          |
| di-n-octyl phthalate      | 10 U          | n-nitroso-di-n-propylamine | 10 U          |
| 1,2-diphenylhydrazine(1)  | 10 U          | n-nitrosodimethylamine     | 10 U          |
| fluoranthene              | 10 U          | n-nitrosodiphenylamine(2)  | 10 U          |
| fluorene                  | 10 U          | phenanthrene               | 10 U          |
| hexachlorobenzene         | 10 U          | pyrene                     | 10 U          |
| hexachlorobutadiene       | 10 U          | 1,2,4-trichlorobenzene     | 10 U          |
| hexachlorocyclopentadiene | 10 U          |                            |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

(1) - Screened for as azobenzene

(2) - Detected as diphenylamine

Job Number: ITPK 51491

## ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-066-03 Lab Sample ID: K3234

| Compound                   | Concentration | Compound              | Concentration |
|----------------------------|---------------|-----------------------|---------------|
| 4-chloro-3-methylphenol    | 10 U          | 2-nitrophenol         | 10 U          |
| 2-chlorophenol             | 10 U          | 4-nitrophenol         | 50 U          |
| 2,4-dichlorophenol         | 10 U          | pentachlorophenol     | <b>50</b> U   |
| 2,4-dimethylphenol         | 10 U          | phenol                | 10 U          |
| 2,4-dinitrophenol          | <b>50</b> U   | 2,4,6-trichlorophenol | 10 U          |
| 2-methyl-4,6-dinitrophenol | <b>50</b> U   |                       | ·             |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Job Number: ITPK 51491

## ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-066-03 Lab Sample ID: K3234

| Tentative Identification (1) | Concentration (2) |
|------------------------------|-------------------|
| 2-pentanone, 4-hydroxy-4-met | 51 AB             |
| benzothiazole                | 9.5               |
| benzothiazole, 2-(methylthio | 6.4               |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- A Suspected aldol condensation product.
- B Analyte was found in the blank as well as the sample.

Job Number: ITPK 51491

#### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-066-04 Lab Sample ID: K3235

| Compound                   | Concentration | Compound                    | Concentration |
|----------------------------|---------------|-----------------------------|---------------|
| acenaphthene               | 1,000 U       | bis(2-chloroisopropyl)ether | 1,000 U       |
| acenaphthylene             | 1,000 U       | bis(2-ethylhexyl)phthalate  | 360 Ј         |
| anthracene                 | 1,000 U       | 4-bromophenyl phenyl ether  | 1,000 U       |
| benzidine                  | 5,000 U       | 2-chloronaphthalene         | 1,000 U       |
| benzo(a)anthracene         | 1,000 U       | 4-chlorophenyl phenyl ether | 1,000 U       |
| benzo(b)fluoranthene       | 1,000 U       | chrysene                    | 1,000 U       |
| benzo(k)fluoranthene       | 1,000 U       | dibenz(a,h)anthracene       | 1,000 U       |
| benzo(a)pyrene             | 1,000 U       | di-n-butylphthalate         | 1,000 U       |
| benzo(g,h,i)perylene       | 1,000 U       | 1,2-dichlorobenzene         | 1,000 U       |
| butylbenzylphthalate       | 1,000 U       | 1,3-dichlorobenzene         | 1,000 U       |
| bis(2-chloroethoxy)methane | 1,000 U       | 1,4-dichlorobenzene         | 1,000 U       |
| bis(2-chloroethyl)ether    | 1,000 U .     |                             |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Job Number: ITPK 51491

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-066-04 Lab Sample ID: K3235

| Compound                  | Concentrati | on       | Compound                   | Concentra | ation |
|---------------------------|-------------|----------|----------------------------|-----------|-------|
| 3,3'dichlorobenzidine     | 2,000       | U        | hexachloroethane           | 1,000     | Ŭ     |
| diethylphthalate          | 1,000       | U        | indeno(1,2,3-cd)pyrene     | 1,000     | Ü     |
| dimethyl phthalate        | 1,000       | <b>U</b> | isophorone                 | 1,000     | U     |
| 2,4-dinitrotoluene        | 1,000       | U        | naphthalene                | 1,000     | U     |
| 2,6-dinitrotoluene        | 1,000       | U        | nitrobenzene               | 1,000     | U     |
| di-n-octyl phthalate      | 1,000       | U        | n-nitroso-di-n-propylamine | 1,000     | Ų     |
| 1,2-diphenylhydrazine(1)  | 1,000       | U        | n-nitrosodimethylamine     | 1,000     | U     |
| fluoranthene              | 1,000       | Ŭ        | n-nitrosodiphenylamine(2)  | 1,000     | U     |
| fluorene                  | 1,000       | U        | phenanthrene               | 1,000     | U     |
| hexachlorobenzene         | 1,000       | U        | pyrene                     | 1,000     | U     |
| hexachlorobutadiene       | 1,000       | U        | 1,2,4-trichlorobenzene     | 1,000     | U     |
| hexachlorocyclopentadiene | 1,000       | U        |                            |           |       |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

<sup>(1) -</sup> Screened for as azobenzene

<sup>(2) -</sup> Detected as diphenylamine

Job Number: ITPK 51491

## ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-066-04 Lab Sample ID: K3235

| Compound                   | Concentration | Compound              | Concentration |
|----------------------------|---------------|-----------------------|---------------|
| 4-chloro-3-methylphenol    | 1,000 U       | 2-nitrophenol         | 1,000 U       |
| 2-chlorophenol             | 1,000 U       | 4-nitrophenol         | 5,000 U       |
| 2,4-dichlorophenol         | 1,000 U       | pentachlorophenol     | 5,000 U       |
| 2,4-dimethylphenol         | 1,000 U       | phenol                | 1,000 U       |
| 2,4-dinitrophenol          | 5,000 U       | 2,4,6-trichlorophenol | 1,000 U       |
| 2-methyl-4,6-dinitrophenol | 5,000 U       |                       | ·             |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Client Project ID: LE Carpenter

Job Number: ITPK 51491

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in  $\mu$ g/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-066-04 Lab Sample ID: K3235

| Tentative Identification (1) | Concentration (2) |
|------------------------------|-------------------|
| unknown                      | 2,500             |
| 9-octadecenoic acid (Z)-, 2, | 1,800             |
| unknown (UNSAT'D?)           | 880               |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITPK 51491

## BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-068-01 Lab Sample ID: K3236

| Compound                   | Concentration | Compound                    | Concentration |
|----------------------------|---------------|-----------------------------|---------------|
| acenaphthene               | 11 U          | bis(2-chloroisopropyl)ether | 11 U          |
| acenaphthylene             | 11 U          | bis(2-ethylhexyl)phthalate  | 11 U          |
| anthracene                 | 11 U          | 4-bromophenyl phenyl ether  | 11 Ŭ          |
| benzidine                  | 53 U          | 2-chloronaphthalene         | 11 U          |
| benzo(a)anthracene         | 11 U          | 4-chlorophenyl phenyl ether | 11 U          |
| benzo(b)fluoranthene       | 11 U          | chrysene                    | 11 U          |
| benzo(k)fluoranthene       | 11 U          | dibenz(a,h)anthracene       | 11 U          |
| benzo(a)pyrene             | 11 · U        | di-n-butylphthalate         | 11 U          |
| benzo(g,h,i)perylene       | 11 U          | 1,2-dichlorobenzene         | 11 U          |
| butylbenzylphthalate       | 3 J           | 1,3-dichlorobenzene         | 11 U          |
| bis(2-chloroethoxy)methane | 11 U          | 1,4-dichlorobenzene         | 11 U          |
| bis(2-chloroethyl)ether    | 11. U         |                             |               |

U Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction:

06/02/92

Date of Analysis: 06/08/92

Indicates an estimated value less than the detection limit.

Job Number: ITPK 51491

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in  $\mu g/\text{liter}$  (ppb)

Sample Matrix: Water

Client Sample ID: 1045-068-01 Lab Sample ID: K3236

| Compound                  | Concentration | Compound                   | Concentration |
|---------------------------|---------------|----------------------------|---------------|
| 3,3'dichlorobenzidine     | 21 U          | hexachloroethane           | 11 U          |
| diethylphthalate          | 11 U          | indeno(1,2,3-cd)pyrene     | 11 U          |
| dimethyl phthalate        | 11 U          | isophorone                 | 11 U          |
| 2,4-dinitrotoluene        | 11 U          | naphthalene                | 11 Ù          |
| 2,6-dinitrotoluene        | 11 U          | nitrobenzene               | 11 U          |
| di-n-octyl phthalate      | 11 U          | n-nitroso-di-n-propylamine | . 11 U        |
| 1,2-diphenylhydrazine(1)  | 11 U          | n-nitrosodimethylamine     | 11 U          |
| fluoranthene              | 11 U          | n-nitrosodiphenylamine(2)  | 11 <b>U</b>   |
| fluorene                  | 11 U          | phenanthrene               | 11 <b>U</b>   |
| hexachlorobenzene         | 11 U          | pyrene                     | 11 <b>U</b>   |
| hexachlorobutadiene       | 11 U          | 1,2,4-trichlorobenzene     | 11 U          |
| hexachlorocyclopentadiene | 11 U          |                            |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

<sup>(1) -</sup> Screened for as azobenzene

<sup>(2) -</sup> Detected as diphenylamine

Job Number: ITPK 51491

# ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-068-01 Lab Sample ID: K3236

| Compound                   | Concentration | Compound              | Concentration |
|----------------------------|---------------|-----------------------|---------------|
| 4-chloro-3-methylphenol    | 11 U          | 2-nitrophenol         | 11 U          |
| 2-chlorophenol             | 11 U          | 4-nitrophenol         | 53 U          |
| 2,4-dichlorophenol         | 11 U          | pentachlorophenol     | / <b>53 U</b> |
| 2,4-dimethylphenol         | 98            | phenol                | 11 U          |
| 2,4-dinitrophenol          | 53 U          | 2,4,6-trichlorophenol | 11 <b>U</b>   |
| 2-methyl-4,6-dinitrophenol | 53 U          |                       |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Client Project ID: LE Carpenter

Job Number: ITPK 51491

## ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-068-01 Lab Sample ID: K3236

| Tentative Identification (1) | Concentration (2) |
|------------------------------|-------------------|
| 2-pentanone, 4-hydroxy-4-met | 60 AB             |
| benzothiazole                | 22                |
| benzene, -ethyl-methyl-      | 31 Y              |
| benzene, -ethyl-methyl-      | 20 Y              |
| benzene, -ethyl-methyl-      | 20 Y              |
| benzene, -ethyl-methyl-      | 14 Y              |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- A Suspected aldol condensation product.
- B Analyte was found in the blank as well as the sample.
- Y Indistinguishable isomer in tentatively identified compounds.

Job Number: ITPK 51491

## BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-068-02 Lab Sample ID: K3237

| Compound                   | Concentration | Compound                    | Concentration |
|----------------------------|---------------|-----------------------------|---------------|
| acenaphthene               | 11 U          | bis(2-chloroisopropyl)ether | 11 U          |
| acenaphthylene             | 11 U          | bis(2-ethylhexyl)phthalate  | 11 U          |
| anthracene                 | 11 U          | 4-bromophenyl phenyl ether  | 11 U          |
| benzidine                  | <b>5</b> 3 U  | 2-chloronaphthalene         | 11 U          |
| benzo(a)anthracene         | 11 U          | 4-chlorophenyl phenyl ether | 11 U          |
| benzo(b)fluoranthene       | 11 <b>U</b>   | chrysene                    | 11 U          |
| benzo(k)fluoranthene       | ii U          | dibenz(a,h)anthracene       | _11 U         |
| benzo(a)pyrene             | 11 U          | di-n-butylphthalate         | 11 U          |
| benzo(g,h,i)perylene       | 11 U          | 1,2-dichlorobenzene         | 11 U          |
| butylbenzylphthalate       | 11 U          | 1,3-dichlorobenzene         | 11 U          |
| bis(2-chloroethoxy)methane | 11 U          | 1,4-dichlorobenzene         | 11 U          |
| bis(2-chloroethyl)ether    | 11 U          |                             |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction:

06/02/92

Date of Analysis:

06/08/92

J - Indicates an estimated value less than the detection limit.

Job Number: ITPK 51491

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-068-02 Lab Sample ID: K3237

| Compound                  | Concentra | tion | Compound                   | Concentra | ation    |
|---------------------------|-----------|------|----------------------------|-----------|----------|
| 3,3'dichlorobenzidine     | 21        | U    | hexachloroethane           | 11        | Ŭ        |
| diethylphthalate          | 11        | U    | indeno(1,2,3-cd)pyrene     | 11        | U        |
| dimethyl phthalate        | 11        | U    | isophorone                 | 11        | U        |
| 2,4-dinitrotoluene        | 11        | U    | naphthalene                | 11        | U        |
| 2,6-dinitrotoluene        | 11        | U    | nitrobenzene               | 11        | Ù        |
| di-n-octyl phthalate      | 11        | U    | n-nitroso-di-n-propylamine | 11        | Ü        |
| 1,2-diphenylhydrazine(1)  | 11        | Ü    | n-nitrosodimethylamine     | 11        | U        |
| fluoranthene              | 11        | U    | n-nitrosodiphenylamine(2)  | 11        | U        |
| fluorene                  | 11        | Ŭ    | phenanthrene               | 11        | <b>ט</b> |
| hexachlorobenzene         | 11        | Ŭ    | pyrene                     | 11        | U        |
| hexachlorobutadiene       | 11        | U    | 1,2,4-trichlorobenzene     | 11.       | U        |
| hexachlorocyclopentadiene | 11        | U    |                            |           |          |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

<sup>(1) -</sup> Screened for as azobenzene

<sup>(2) -</sup> Detected as diphenylamine

Job Number: ITPK 51491

# ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-068-02 Lab Sample ID: K3237

| Compound                   | Concentration | Compound              | Concentration |
|----------------------------|---------------|-----------------------|---------------|
| 4-chloro-3-methylphenol    | 11 U          | 2-nitrophenol         | 11 Ü          |
| 2-chlorophenol             | 11 Ú          | 4-nitrophenol         | 53 U          |
| 2,4-dichlorophenol         | 11 U          | pentachlorophenol     | 53 U          |
| 2,4-dimethylphenol         | 11 U          | phenol                | 11 U          |
| 2,4-dinitrophenol          | 53 U          | 2,4,6-trichlorophenol | 11 U          |
| 2-methyl-4,6-dinitrophenol | 53 U          |                       |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Client Project ID: LE Carpenter

Job Number: ITPK 51491

## ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-068-02 Lab Sample ID: K3237

| Tentative Identification (1) | Concentration (2) |
|------------------------------|-------------------|
| 2-pentanone, 4-hydroxy-4-met | 46 AB             |
| benzothiazole                | 22                |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- A Suspected aldol condensation product.
- B Analyte was found in the blank as well as the sample.

Job Number: ITPK 51491

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in  $\mu g/liter$  (ppb)

Sample Matrix: Water

Client Sample ID: 1045-068-03 Lab Sample ID: K3238

| Compound                   | Concentration | Compound                    | Concentration |
|----------------------------|---------------|-----------------------------|---------------|
| acenaphthene               | 11 U          | bis(2-chloroisopropyl)ether | 11 U          |
| acenaphthylene             | 11 U          | bis(2-ethylhexyl)phthalate  | 11 U          |
| anthracene                 | <b>11 U</b>   | 4-bromophenyl phenyl ether  | 11 U          |
| benzidine                  | <b>5</b> 3 U  | 2-chloronaphthalene         | 11 U          |
| benzo(a)anthracene         | 11 U          | 4-chlorophenyl phenyl ether | 11 U          |
| benzo(b)fluoranthene       | 11 U          | chrysene                    | 11 U          |
| benzo(k)fluoranthene       | 11 U          | dibenz(a,h)anthracene       | 11 U          |
| benzo(a)pyrene             | 11. U         | di-n-butylphthalate         | 11 U          |
| benzo(g,h,i)perylene       | 11 U          | 1,2-dichlorobenzene         | 11 U          |
| butylbenzylphthalate       | 2 J           | 1,3-dichlorobenzene         | 11 U          |
| bis(2-chloroethoxy)methane | 11 U          | 1,4-dichlorobenzene         | 11 U          |
| bis(2-chloroethyl)ether    | 11 U          |                             |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Job Number: ITPK 51491

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-068-03 Lab Sample ID: K3238

| Compound                  | Concentration | Compound                   | Concentration |
|---------------------------|---------------|----------------------------|---------------|
| 3,3'dichlorobenzidine     | 21 U          | hexachloroethane           | 11 Ŭ          |
| diethylphthalate          | 11 U          | indeno(1,2,3-cd)pyrene     | 11 U          |
| dimethyl phthalate        | 11. U         | isophorone                 | 11 U          |
| 2,4-dinitrotoluene        | 11 Ū          | naphthalene                | 11 · U        |
| 2,6-dinitrotoluene        | 11 U          | nitrobenzene               | 11 U          |
| di-n-octyl phthalate      | 11 U          | n-nitroso-di-n-propylamine | 11 U          |
| 1,2-diphenylhydrazine(1)  | 11 U          | n-nitrosodimethylamine     | 11 U          |
| fluoranthene              | 11 <b>V</b>   | n-nitrosodiphenylamine(2)  | 11 U          |
| fluorene                  | 11 Ü          | phenanthrene               | 11 U          |
| hexachlorobenzene         | 11 Ü          | pyrene                     | 11 <b>U</b>   |
| hexachloróbutadiene       | 11 U          | 1,2,4-trichlorobenzene     | 11 <b>U</b>   |
| hexachlorocyclopentadiene | 11 U          |                            |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

- Indicates an estimated value less than the detection limit.

(1) - Screened for as azobenzene

(2) - Detected as diphenylamine

Job Number: ITPK 51491

#### ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-068-03 Lab Sample ID: K3238

| Compound                   | Concentration | Compound              | Concentration |
|----------------------------|---------------|-----------------------|---------------|
| 4-chloro-3-methylphenol    | 11 U          | 2-nitrophenol         | 11 U          |
| 2-chlorophenol             | 11 U          | 4-nitrophenol         | <b>53</b> Û   |
| 2,4-dichlorophenol         | 11 U          | pentachlorophenol     | 53 U          |
| 2,4-dimethylphenol         | 11 U          | phenol                | 11 U          |
| 2,4-dinitrophenol          | 53 U          | 2,4,6-trichlorophenol | 11 Ü          |
| 2-methyl-4,6-dinitrophenol | 53 U          | ,                     |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Client Project ID: LE Carpenter

Job Number: ITPK 51491

# ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in  $\mu g$ /liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-068-03 Lab Sample ID: K3238

| Tentative Identification (1) | Concentration (2) |
|------------------------------|-------------------|
| benzothiazole                | 33                |
| 2-pentanone, 4-hydroxy-4-met | 12 AB             |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- A Suspected aldol condensation product.
- B Analyte was found in the blank as well as the sample.

Job Number: ITPK 51491

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-068-04 Lab Sample ID: K3239

| Compound                   | Concentration | Compound                    | Concentration |
|----------------------------|---------------|-----------------------------|---------------|
| acenaphthene               | 13 U          | bis(2-chloroisopropyl)ether | 13 U          |
| acenaphthylene             | 13 U          | bis(2-ethylhexyl)phthalate  | 13 Ú          |
| anthracene                 | 13 U          | 4-bromophenyl phenyl ether  | 13 U          |
| benzidine                  | 64 U          | 2-chloronaphthalene         | 13 U          |
| benzo(a)anthracene         | 13 U          | 4-chlorophenyl phenyl ether | 13 U          |
| benzo(b)fluoranthene       | 13 U          | chrysene                    | 13 U          |
| benzo(k)fluoranthene       | 13 U          | dibenz(a,h)anthracene       | 13 U          |
| benzo(a)pyrene             | 13 U          | di-n-butylphthalate         | 13 U          |
| benzo(g,h,i)perylene       | 13 U          | 1,2-dichlorobenzene         | 13 U          |
| butylbenzylphthalate       | 13 U          | 1,3-dichlorobenzene         | 13 U          |
| bis(2-chloroethoxy)methane | 13 U          | 1,4-dichlorobenzene         | 13 Ü          |
| bis(2-chloroethyl)ether    | 13 U          |                             |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction:

06/02/92

Date of Analysis:

06/08/92

J - Indicates an estimated value less than the detection limit.

Job Number: ITPK 51491

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-068-04 Lab Sample ID: K3239

| Compound                  | Concentration   | Compound                   | Concentration |
|---------------------------|-----------------|----------------------------|---------------|
| 3,3'dichlorobenzidine     | 26 Ü            | hexachloroethane           | 13 U          |
| diethylphthalate          | 13 U            | indeno(1,2,3-cd)pyrene     | 13 U          |
| dimethyl phthalate        | 13 U            | isophorone                 | 13 U          |
| 2,4-dinitrotoluene        | 13 U            | naphthalene                | 13 U          |
| 2,6-dinitrotoluene        | 13 U            | nitrobenzene               | 13 U          |
| di-n-octyl phthalate      | 13 Ü            | n-nitroso-di-n-propylamine | 13 U          |
| 1,2-diphenylhydrazine(1)  | 13 U            | n-nitrosodimethylamine     | 13 U          |
| fluoranthene              | 13 U            | n-nitrosodiphenylamine(2)  | 13 U          |
| fluorene                  | 13 U            | phenanthrene               | 13 U          |
| hexachlorobenzene         | 13 Ü            | pyrene                     | 13 U          |
| hexachlorobutadiene       | 13 Ü            | 1,2,4-trichlorobenzene     | 13 U          |
| hexachlorocyclopentadiene | / <b>13 U</b> F |                            |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

<sup>(1) -</sup> Screened for as azobenzene

<sup>(2) -</sup> Detected as diphenylamine

Job Number: ITPK 51491

## ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-068-04 Lab Sample ID: K3239

| Compound                   | Concen | tration | Compound              | Concentration |   |
|----------------------------|--------|---------|-----------------------|---------------|---|
| 4-chloro-3-methylphenol    | 13     | U       | 2-nitrophenol         | 13            | U |
| 2-chlorophenol             | 13     | Ų       | 4-nitrophenol         | 64            | U |
| 2,4-dichlorophenol         | 13     | U       | pentachlorophenol     | 64            | U |
| 2,4-dimethylphenol         | 13     | U       | phenol                | 13            | Ü |
| 2,4-dinitrophenol          | 64     | U ;     | 2,4,6-trichlorophenol | 13            | U |
| 2-methyl-4,6-dinitrophenol | 64     | U       |                       |               |   |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 0

06/02/92

Date of Analysis:

06/08/92

J - Indicates an estimated value less than the detection limit.

Client Project ID: LE Carpenter

Job Number: ITPK 51491

# ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-068-04 Lab Sample ID: K3239

Tentative Identification (1)

Concentration (2)

None Detected

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITPK 51491

## BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank

Lab Sample ID: H0679

| Compound                   | Concentration | Compound                    | Concentration |
|----------------------------|---------------|-----------------------------|---------------|
| acenaphthene               | 10 U          | bis(2-chloroisopropyl)ether | 10 U          |
| acenaphthylene             | 10 U          | bis(2-ethylhexyl)phthalate  | 10 U          |
| anthracene                 | 10 U          | 4-bromophenyl phenyl ether  | 10 U          |
| benzidine                  | 50 U          | 2-chloronaphthalene         | 10 Ŭ          |
| benzo(a)anthracene         | 10 U          | 4-chlorophenyl phenyl ether | 10 U          |
| benzo(b)fluoranthene       | 10 U          | chrysene                    | 10 U          |
| benzo(k)fluoranthene       | 10 U          | dibenz(a,h)anthracene       | 10 U          |
| benzo(a)pyrene             | 10 U          | di-n-butylphthalate         | 10 Ü          |
| benzo(g,h,i)perylene       | 10 U          | 1,2-dichlorobenzene         | 10 U          |
| butylbenzylphthalate       | 10 U          | 1,3-dichlorobenzene         | 10 U          |
| bis(2-chloroethoxy)methane | 10 U          | 1,4-dichlorobenzene         | 10 U          |
| bis(2-chloroethyl)ether    | 10 U          |                             |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Job Number: ITPK 51491

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank Lab Sample ID: H0679

| Compound                  | Concentration | Compound                   | Concentration |
|---------------------------|---------------|----------------------------|---------------|
| 3,3'dichlorobenzidine     | 20 U          | hexachloroethane           | 10 U          |
| diethylphthalate          | 10 U          | indeno(1,2,3-cd)pyrene     | 10 U          |
| dimethyl phthalate        | 10 U          | isophorone                 | 10 U          |
| 2,4-dinitrotoluene        | 10 Ų          | naphthalene                | 10 U          |
| 2,6-dinitrotoluene        | . 10 U        | nitrobenzene               | 10 U          |
| di-n-octyl phthalate      | 10 U          | n-nitroso-di-n-propylamine | 10 U          |
| 1,2-diphenylhydrazine(1)  | 10 U          | n-nitrosodimethylamine     | 10 U          |
| fluoranthene              | , 10 Ŭ        | n-nitrosodiphenylamine(2)  | 10 U          |
| fluorene                  | 10 U          | phenanthrene               | 10 U          |
| hexachlorobenzene         | 10 U          | pyrene                     | 10 Ü          |
| hexachlorobutadiene       | 10 U          | 1,2,4-trichlorobenzene     | 10 U          |
| hexachlorocyclopentadiene | 10 U          |                            |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 0
Date of Analysis: 0

06/02/92

06/08/92

J - Indicates an estimated value less than the detection limit.

<sup>(1) -</sup> Screened for as azobenzene

<sup>(2) -</sup> Detected as diphenylamine

Job Number: ITPK 51491

#### ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank

Lab Sample ID: H0679

| Compound                   | Concentration | Compound              | Concentration |  |
|----------------------------|---------------|-----------------------|---------------|--|
| 4-chloro-3-methylphenol    | 10 U          | 2-nitrophenol         | 10 U          |  |
| 2-chlorophenol             | 10 U          | 4-nitrophenol         | 50 U          |  |
| 2,4-dichlorophenol         | 10 U          | pentachlorophenol     | 50 U          |  |
| 2,4-dimethylphenol         | 10 U          | phenol                | 10 U          |  |
| 2,4-dinitrophenol          | 50 U          | 2,4,6-trichlorophenol | 10 U          |  |
| 2-methyl-4,6-dinitrophenol | 50 U          |                       |               |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Indicates an estimated value less than the detection limit.

Client Project ID: LE Carpenter

Job Number: ITPK 51491

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank Lab Sample ID: H0679

Tentative Identification (1)

Concentration (2)

2-pentanone, 4-hydroxy-4-methyl-

32 A

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- A Suspected aldol condensation product.

Job Number: ITPK 51491

## WATER SURROGATE PERCENT RECOVERY SUMMARY

## **SEMI-VOLATILE**

| Client Sample ID | Nitro-<br>Benzene-D5<br>(35-114%)* | 2-Fluoro-<br>Biphenyl<br>(43-116%)* | Terphenyl-<br>D14<br>(33-141%)* | Phenol-D5<br>(10-94%)* | 2-Fluoro-<br>Phenol<br>(21-100%)* | 2,4,6-<br>Tribromo-<br>Phenol<br>(10-123%)* |
|------------------|------------------------------------|-------------------------------------|---------------------------------|------------------------|-----------------------------------|---|
| 1045-066-01      | D                                  | D                                   | D                               | D                      | D                                 | D   |
| 1045-066-02      | 91                                 | 86                                  | 86                              | 40                     | 53                                | 62  |
| 1045-066-03      | 89                                 | 87                                  | 76                              | 43                     | 53                                | 58  |
| 1045-066-04      | D                                  | D                                   | D                               | D                      | D                                 | D   |
| 1045-068-01      | 93                                 | 84                                  | 76                              | 43                     | 52                                | 61  |
| 1045-068-02      | 89                                 | 81                                  | 69                              | 41                     | 47                                | <b>55</b>                                   |
| 1045-068-03      | 81                                 | 77                                  | 70                              | 40                     | 50                                | 59  |
| 1045-068-04      | 83                                 | 81                                  | 74                              | 44                     | 50                                | 58  |
| Method Blank     | 93                                 | 88                                  | 85                              | - 41                   | .55                               | 64  |

<sup>\*</sup>Values in parenthesis represent QC limits.

D - Surrogates diluted out

Client Project ID: LE Carpenter

Job Number: ITPK 51499

## **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in µg/liter (ppb) Sample Matrix: Water

Client Sample ID: 1045-069-01 Lab Sample ID: K3274

| Compound                 | Concentration | Compound                  | Concentration |
|--------------------------|---------------|---------------------------|---------------|
| acrolein                 | 1,000 U       | 1,1-dichloroethene        | 500 U         |
| acrylonitrile            | 1,000 U       | cis-1,2-dichloroethene    | 500 U         |
| benzene                  | 500 U         | trans-1,2-dichloroethene  | 500 Ü         |
| bromodichloromethane     | <b>500</b> U  | 1,2-dichloropropane       | 500 U         |
| bromoform                | 500 U         | cis-1,3-dichloropropene   | 500 U         |
| bromomethane             | 1,000 U       | trans-1,3-dichloropropene | 500 U         |
| carbon tetrachloride     | 500 U         | ethyl benzene             | <b>500</b> U  |
| chlorobenzene            | 500 U         | methylene chloride        | 1,800 B       |
| chloroethane             | 1,000 U       | 1,1,2,2-tetrachloroethane | 500 U         |
| 2-chloroethylvinyl ether | 1,000 U       | tetrachloroethene         | 500 U         |
| chloroform               | <b>500</b> U  | toluene                   | 500 U         |
| chloromethane            | 1,000 U       | 1,1,1-trichloroethane     | 500 U         |
| dibromochloromethane     | 500 U         | 1,1,2-trichloroethane     | <b>500</b> U  |
| 1,1-dichloroethane       | 500 U         | trichloroethene           | 500 U         |
| 1,2-dichloroethane       | 500 U         | trichlorofluoromethane    | 500 U         |
|                          |               | vinyl chloride            | 1,000 U       |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/05/92

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

Client Project ID: LE Carpenter

Job Number: ITPK 51499

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in  $\mu g/liter$  (ppb)

Sample Matrix: Water

Client Sample ID: 1045-069-01 Lab Sample ID: K3274

| Tentative Identification (1) | Concentration (2) |
|------------------------------|-------------------|
| benzene, dimethyl-           | 5,700 Y           |
| benzene, 1,2-dimethyl-       | 2,900             |
| unknown (substituted alkane) | .520              |
| unknown (alcohol?)           | 740               |
| decane                       | <b>940</b>        |
| undecane                     | 1,400             |

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Y - Indistinguishable isomer in tentatively identified compounds.

Job Number: ITPK 51499

# **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g/liter$  (ppb) Sample Matrix: Water

Client Sample ID: 1045-069-02 Lab Sample ID: K3275

| Compound                 | Concent | ration | Compound                  | Concent | ration |
|--------------------------|---------|--------|---------------------------|---------|--------|
| acrolein                 | 10      | Ü      | 1,1-dichloroethene        | 5       | U      |
| acrylonitrile            | 10      | U      | cis-1,2-dichloroethene    | 5       | U      |
| benzene                  | 6       |        | trans-1,2-dichloroethene  | 5       | U      |
| bromodichloromethane     | 5       | U      | 1,2-dichloropropane       | 5       | U      |
| bromoform                | 5       | U      | cis-1,3-dichloropropene   | 5       | U      |
| bromomethane             | 10      | U      | trans-1,3-dichloropropene | 5       | U      |
| carbon tetrachloride     | 5       | U      | ethyl benzene             | 3       | J      |
| chlorobenzene            | 5       | U      | methylene chloride        | 410     | DB     |
| chloroethane             | 10      | U      | 1,1,2,2-tetrachloroethane | 5       | Ü      |
| 2-chloroethylvinyl ether | 10      | U      | tetrachloroethene         | 5       | U      |
| chloroform               | 6       |        | toluene                   | 14      |        |
| chloromethane            | 10      | U      | 1,1,1-trichloroethane     | 5       | U      |
| dibromochloromethane     | 5       | Ü      | 1,1,2-trichloroethane     | 5       | U      |
| 1,1-dichloroethane       | 5       | U      | trichloroethene           | 5       | U      |
| 1,2-dichloroethane       | 5       | U      | trichlorofluoromethane    | 5       | U      |
|                          |         | ĸ      | vinyl chloride            | 10      | U      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/08/92

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

D - Compound analyzed at a secondary dilution factor.

Client Project ID: LE Carpenter Job Number: ITPK 51499

# ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-069-02 Lab Sample ID: K3275

| Tentative Identification (1) | Concentration (2) |  |  |
|------------------------------|-------------------|--|--|
| benzene, dimethyl-           | 50 Y              |  |  |
| benzene, 1,2-dimethyl-       | 9.2               |  |  |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- Y Indistinguishable isomer in tentatively identified compounds.

Job Number: ITPK 51499

# **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: 1045-069-03 Lab Sample ID: K3276

| Compound                 | Concentrat          | <u>tion</u> | Compound                  | Concent  | ration |
|--------------------------|---------------------|-------------|---------------------------|----------|--------|
| acrolein                 | 10 T                | U           | 1,1-dichloroethene        | 5        | U      |
| acrylonitrile            | 10 T                | U           | cis-1,2-dichloroethene    | 5        | U      |
| benzene                  | 5 T                 | Ŭ           | trans-1,2-dichloroethene  | <b>5</b> | U      |
| bromodichloromethane     | 2 J                 | 7           | 1,2-dichloropropane       | 5        | U      |
| bromoform                | <b>5</b> ∋7         | ט           | cis-1,3-dichloropropene   | 5        | U      |
| bromomethane             | 10 T                | U           | trans-1,3-dichloropropene | 5        | U      |
| carbon tetrachloride     | <b>5</b> T          | Ŭ           | ethyl benzene             | 2        | J      |
| chlorobenzene            | .5 ,                | ט           | methylene chloride        | 4,800    | DB     |
| chloroethane             | 10 U                | Ŭ.          | 1,1,2,2-tetrachloroethane | 5        | U      |
| 2-chloroethylvinyl ether | 10 T                | ט           | tetrachloroethene         | 5        | U      |
| chloroform               | 22                  |             | toluene                   | 5        | U      |
| chloromethane            | 10 T                | U           | 1,1,1-trichloroethane     | 5        | U      |
| dibromochloromethane     | <b>5</b> , <b>t</b> | ט           | 1,1,2-trichloroethane     | 5        | U      |
| 1,1-dichloroethane       | 5 T                 | U           | trichloroethene           | 5        | U      |
| 1,2-dichloroethane       | 5 T                 | U           | trichlorofluoromethane    | 5        | U      |
|                          |                     |             | vinyl chloride            | 10       | U      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/05/92

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

D - Compound analyzed at a secondary dilution factor.

Job Number: ITPK 51499

Client Project ID: LE Carpenter

# ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-069-03 Lab Sample ID: K3276

| Tentative Identification (1) | Concentration (2) |
|------------------------------|-------------------|
| benzene, dimethyl-           | 23 Y              |
| benzene, 1,2-dimethyl-       | 7.7               |

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- Y Indistinguishable isomer in tentatively identified compounds.

Client Project ID: LE Carpenter

Job Number: ITPK 51499

# **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: 1045-069-04 Lab Sample ID: K3277

| Compound                 | Concentration | Compound                  | Concentration |
|--------------------------|---------------|---------------------------|---------------|
| acrolein                 | 50 U          | 1,1-dichloroethene        | 25 U          |
| acrylonitrile            | 50 U          | cis-1,2-dichloroethene    | <b>25</b> U   |
| benzene                  | 13 J          | trans-1,2-dichloroethene  | 25 Ü          |
| bromodichloromethane     | 25 U          | 1,2-dichloropropane       | 25 U          |
| bromoform                | 25 U          | cis-1,3-dichloropropene   | 25 U          |
| bromomethane             | <b>50</b> U   | trans-1,3-dichloropropene | 25 U          |
| carbon tetrachloride     | 25 U          | ethyl benzene             | 25 U          |
| chlorobenzene            | 25 Ù          | methylene chloride        | 68 B          |
| chloroethane             | 50 U          | 1,1,2,2-tetrachloroethane | 25 U          |
| 2-chloroethylvinyl ether | 50 U          | tetrachloroethene         | <b>25</b> U   |
| chloroform               | 21 J          | tõluene                   | 20 J          |
| chloromethane            | 50 U          | 1,1,1-trichloroethane     | 25 U          |
| dibromochloromethane     | 25 U          | 1,1,2-trichloroethane     | 25 U          |
| 1,1-dichloroethane       | 25 U          | trichloroethene           | · 25 U        |
| 1,2-dichloroethane       | 25 Ü          | trichlorofluoromethane    | 25 U          |
|                          |               | vinyl chloride            | <b>50</b> U   |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/05/92

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

Job Number: ITPK 51499

Client Project ID: LE Carpenter

### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-069-04 Lab Sample ID: K3277

| Tentative Identification (1) | Concentration (2) |
|------------------------------|-------------------|
| pentane                      | 66                |
| hexane                       | 68                |
| octane                       | <b>56</b>         |
| benzene, dimethyl-           | 41 Y              |

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

- Indistinguishable isomer in tentatively identified compounds.

Job Number: ITPK 51499

### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: Method Blank 1

Lab Sample ID: WB0605

| Compound                 | Concentration | Compound                  | Concentration |
|--------------------------|---------------|---------------------------|---------------|
| acrolein                 | 10 Ù          | 1,1-dichloroethene        | 5 U           |
| acrylonitrile            | 10 "U         | cis-1,2-dichloroethene    | 5 U           |
| benzene                  | 5 U           | trans-1,2-dichloroethene  | 5 U           |
| bromodichloromethane     | 5 . U         | 1,2-dichloropropane       | 5 U           |
| bromoform                | 5 U           | cis-1,3-dichloropropene   | 5 U           |
| bromomethane             | 10 U          | trans-1,3-dichloropropene | 5 U           |
| carbon tetrachloride     | <b>5</b> U    | ethyl benzene             | 5 U           |
| chlorobenzene            | 5 U           | methylene chloride        | 2 J           |
| chloroethane             | 10 U          | 1,1,2,2-tetrachloroethane | 5 U           |
| 2-chloroethylvinyl ether | 10 U          | tetrachloroethene         | 5 U           |
| chloroform               | 5 U           | toluene                   | 5 U           |
| chloromethane            | 10 U          | 1,1,1-trichloroethane     | 5 Ù           |
| dibromochloromethane     | 5 U           | 1,1,2-trichloroethane     | 5 U           |
| 1,1-dichloroethane       | 5 U           | trichloroethene           | 5 U           |
| 1,2-dichloroethane       | 5 U           | trichlorofluoromethane    | 5 Ù           |
|                          |               | vinyl chloride            | 10 U          |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/05/92

This method blank applies to the following samples: 1045-069-01, 1045-069-02 DL, 1045-069-03, 1045-069-03 DL.

DL - Dilution

J - Indicates an estimated value less than the detection limit.

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

Job Number: ITPK 51499

### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 1

Lab Sample ID: WB0605

Tentative Identification (1)

Concentration (2)

None Detected

Remarks:

(1) Identification is based on computer search of the NIST Library.

(2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Client Project ID: LE Carpenter

Job Number: ITPK 51499

# **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in  $\mu g$ /liter (ppb) Sample Matrix: Water

Client Sample ID: Method Blank 2 Lab Sample ID: WB06052

| Compound                 | Concentration |          | Compound                  | Concent | ration       |
|--------------------------|---------------|----------|---------------------------|---------|--------------|
| acrolein                 | 10            | U        | 1,1-dichloroethene        | 5       | U            |
| acrylonitrile            | 10            | <b>U</b> | cis-1,2-dichloroethene    | 5       | · <b>U</b> . |
| benzene                  | 5             | Ū        | trans-1,2-dichloroethene  | 5       | U            |
| bromodichloromethane     | 5             | U        | 1,2-dichloropropane       | 5       | U            |
| bromoform                | 5             | Ū        | cis-1,3-dichloropropene   | 5       | U            |
| bromomethane             | 10            | U        | trans-1,3-dichloropropene | 5       | Ų            |
| carbon tetrachloride     | 5             | Ù        | ethyl benzene             | 5       | U            |
| chlorobenzene            | 5             | U        | methylene chloride        | 2       | J            |
| chloroethane             | 10            | U        | 1,1,2,2-tetrachloroethane | 5       | U            |
| 2-chloroethylvinyl ether | 10            | U        | tetrachloroethene         | 5       | U            |
| chloroform               | 5             | U        | toluene                   | 5       | U            |
| chloromethane            | 10            | U        | 1,1,1-trichloroethane     | 5       | U            |
| dibromochloromethane     | 5             | U        | 1,1,2-trichloroethane     | 5       | U            |
| 1,1-dichloroethane       | 5             | U        | trichloroethene           | 5       | U            |
| 1,2-dichloroethane       | 5             | U        | trichlorofluoromethane    | 5       | U            |
| •                        |               | 1 11     | vinyl chloride            | 10      | Ų            |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/05/92

This method blank applies to the following sample: 1045-069-04.

J - Indicates an estimated value less than the detection limit.

Job Number: ITPK 51499

Client Project ID: LE Carpenter

# ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 2 Lab Sample ID: WB06052

Tentative Identification (1)

Concentration (2)

None Detected

Remarks:

(1) Identification is based on computer search of the NIST Library.

(2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITPK 51499

# **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in µg/liter (ppb) Sample Matrix: Water

Client Sample ID: Method Blank 3

Lab Sample ID: WB0608

| Compound                 | Concent | ration | Compound                  | Concent | ration     |
|--------------------------|---------|--------|---------------------------|---------|------------|
| acrolein                 | 10      | U      | 1,1-dichloroethene        | 5       | U          |
| acrylonitrile            | 10      | U      | cis-1,2-dichloroethene    | 5       | U          |
| benzene                  | 5       | U      | trans-1,2-dichloroethene  | 5       | Ų          |
| bromodichloromethane     | 5       | Ü      | 1,2-dichloropropane       | 5       | U,         |
| bromoform                | 5       | Ū      | cis-1,3-dichloropropene   | 5       | Ü          |
| bromomethane             | 10      | Ü      | trans-1,3-dichloropropene | 5       | · <b>U</b> |
| carbon tetrachloride     | 5       | U      | ethyl benzene             | 5       | U          |
| chlorobenzene            | 5       | Ŭ      | methylene chloride        | 2       | J          |
| chloroethane             | 10      | U      | 1,1,2,2-tetrachloroethane | 5       | Ü          |
| 2-chloroethylvinyl ether | 10      | Ü      | tetrachloroethene         | . 5     | U          |
| chloroform               | 5       | U      | toluene                   | 5       | U          |
| chloromethane            | 10      | Ŭ      | 1,1,1-trichloroethane     | 5       | U          |
| dibromochloromethane     | 5       | U      | 1,1,2-trichloroethane     | 5       | U          |
| 1,1-dichloroethane       | 5       | U      | trichloroethene           | . 5     | U          |
| 1,2-dichloroethane       | 5       | U      | trichlorofluoromethane    | 5       | U          |
|                          |         |        | vinyl chloride            | 10      | U          |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/08/92

This method blank applies to the following sample: 1045-069-02.

J - Indicates an estimated value less than the detection limit.

Job Number: ITPK 51499

KIOK VIIIEL, IN

Client Project ID: LE Carpenter

# ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 3

Lab Sample ID: WB0608

Tentative Identification (1)

Concentration (2)

None Detected

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITPK 51499

# WATER SURROGATE PERCENT RECOVERY SUMMARY

|                  | <del></del>           | VOLATILE          |                                     |
|------------------|-----------------------|-------------------|-------------------------------------|
| Client Sample ID | Toluene-D8 (88-110%)* | BFB<br>(86-115%)* | 1,2 Dichloroethane-D4<br>(76-114%)* |
| 1045-069-01      | 98                    | 95                | 98                                  |
| 1045-069-02      | 103                   | 99                | 100                                 |
| 1045-069-02 DL   | 96                    | 95                | 100                                 |
| 1045-069-03      | <b>95</b>             | 93                | 91                                  |
| 1045-069-03 DL   | 103                   | 102               | 104                                 |
| 1045-069-04      | 109                   | 90                | 94                                  |
| Method Blank 1   | 97                    | 93                | 96                                  |
| Method Blank 2   | 101                   | 97                | 98                                  |
| Method Blank 3   | 102                   | 96                | 97                                  |

<sup>\*</sup>Values in parenthesis represent QC limits.

DL - Dilution \

Job Number: ITPK 51499

# **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in total  $\mu g$ Sample Matrix: Water

Client Sample ID: Method Blank 1 Lab Sample ID: WB06192

| Compound                 | Concentratio  | n Compound                | Concent | ration     |
|--------------------------|---------------|---------------------------|---------|------------|
| acrolein                 | 10 U          | 1,1-dichloroethene        | 5       | U          |
| acrylonitrile            | 10 U          | cis-1,2-dichloroethene    | 5       | U          |
| benzene                  | 5 U           | trans-1,2-dichloroethene  | 5       | U          |
| bromodichloromethane     | 5 U           | 1,2-dichloropropane       | 5       | U          |
| bromoform                | 5 U           | cis-1,3-dichloropropene   | 5       | U          |
| bromomethane             | 10 U          | trans-1,3-dichloropropene | 5       | U          |
| carbon tetrachloride     | . <b>5</b> 'U | ethyl benzene             | 5       | U          |
| chlorobenzene            | 5 Ŭ           | methylene chloride        | 2       | J          |
| chloroethane             | 10 U          | 1,1,2,2-tetrachloroethane | 5       | U          |
| 2-chloroethylvinyl ether | 10 U          | tetrachloroethene         | . 5     | U          |
| chloroform               | 5 U           | toluene                   | 5       | <b>U</b> . |
| chloromethane            | 10 U          | 1,1,1-trichloroethane     | 5       | U          |
| dibromochloromethane     | 5 U           | 1,1,2-trichloroethane     | 5       | U          |
| 1,1-dichloroethane       | 5 U           | trichloroethene           | 5       | Ù          |
| 1,2-dichloroethane       | 5 Ü           | trichlorofluoromethane    | 5       | U          |
|                          |               | vinyl chloride            | 10      | U          |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/19/92

This method blank applies to the following samples: 1045-069-05B, 1045-069-05T, 1045-069-06T, 1045-069-07T, 1045-069-07T DL, 1045-069-07T DL2, 1045-069-08T.

DL - Dilution.

J - Indicates an estimated value less than the detection limit.

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

Job Number: ITPK 51499

### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in Total µg

Sample Matrix: Water

Client Sample ID: Method Blank 1 Lab Sample ID: WB06192

Tentative Identification (1)

None Detected

Concentration (2)

Remarks:

(1) Identification is based on computer search of the NIST Library.

(2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITPK 51499

# **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in Total  $\mu g$ Sample Matrix: Water

Client Sample ID: Method Blank 2 Lab Sample ID: WB0622R2

| Compound                 | Concentration   | Compound                  |     | Concent | ration | ļ |
|--------------------------|-----------------|---------------------------|-----|---------|--------|---|
| acrolein                 | 10 <b>U</b>     | 1,1-dichloroethene        |     | 5       | U      |   |
| acrylonitrile            | 10 U            | cis-1,2-dichloroethene    |     | 5       | U      |   |
| benzene                  | <b>5</b> U      | trans-1,2-dichloroethene  |     | 5       | U      |   |
| bromodichloromethane     | 5 : U           | 1,2-dichloropropane       |     | 5       | U      |   |
| bromoform                | 5 U             | cis-1,3-dichloropropene   | •   | 5       | ·U     |   |
| bromomethane             | 10 U            | trans-1,3-dichloropropene |     | . 5     | U      |   |
| carbon tetrachloride     | 5 U             | ethyl benzene             |     | 5       | U      |   |
| chlorobenzene            | <b>5 U</b>      | methylene chloride        |     | 2       | J      |   |
| chloroethane             | 10 U            | 1,1,2,2-tetrachloroethane |     | 5       | U,     |   |
| 2-chloroethylvinyl ether | 10 U            | tetrachloroethene         |     | 5       | U      |   |
| chloroform               | <b>5</b> U      | toluene                   |     | . 5     | U      |   |
| chloromethane            | 10 U            | 1,1,1-trichloroethane     |     | 5       | U      |   |
| dibromochloromethane     | <b>5</b> U      | 1,1,2-trichloroethane     | · · | 5       | Ü      |   |
| 1,1-dichloroethane       | <b>5</b> U      | trichloroethene           |     | 5       | U      |   |
| 1,2-dichloroethane       | 5 U             | trichlorofluoromethane    |     | 5       | U      | • |
|                          | . 9<br>1 9<br>1 | vinyl chloride            |     | 10      | U      |   |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/22/92

This method blank applies to the following samples: 1045-069-05B DL, 1045-069-05B DL2, 1045-069-05M, 1045-069-05M DL.

DL - Dilution.

J - Indicates an estimated value less than the detection limit.

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

Job Number: ITPK 51499

### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in Total µg

Sample Matrix: Water

Client Sample ID: Method Blank 2 Lab Sample ID: WB0622R2

Tentative Identification (1)

Concentration (2)

None Detected

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITPK 51499

### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in Total  $\mu g$ Sample Matrix: Water

Client Sample ID: Method Blank 3 Lab Sample ID: WB06224

| Compound                 | Concentration | Compound                  | Concent  | ration |
|--------------------------|---------------|---------------------------|----------|--------|
| acrolein                 | 10 U          | 1,1-dichloroethene        | 5        | U      |
| acrylonitrile            | 10 Ŭ          | cis-1,2-dichloroethene    | 5        | U      |
| benzene                  | 5 ⊎ <b>U</b>  | trans-1,2-dichloroethene  | 5        | U ·    |
| bromodichloromethane     | 5 Ü           | 1,2-dichloropropane       | 5        | U      |
| bromoform                | 5 U           | cis-1,3-dichloropropene   | 5        | U      |
| bromomethane             | 10 U          | trans-1,3-dichloropropene | 5        | U      |
| carbon tetrachloride     | 5 U           | ethyl benzene             | 5        | U      |
| chlorobenzene            | 5 U           | methylene chloride        | 2        | J      |
| chloroethane             | 10 U          | 1,1,2,2-tetrachloroethane | <b>5</b> | U      |
| 2-chloroethylvinyl ether | 10 U          | tetrachloroethene         | 5        | U      |
| chloroform               | 5 U           | toluene                   | 5        | U .    |
| chloromethane            | 10 U          | 1,1,1-trichloroethane     | 5        | U      |
| dibromochloromethane     | 5 U           | 1,1,2-trichloroethane     | 5        | U      |
| 1,1-dichloroethane       | 5 U           | trichloroethene           | 5        | U      |
| 1,2-dichloroethane       | 5 U           | trichlorofluoromethane    | 5        | U      |
|                          |               | vinyl chloride            | 10       | U      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/22/92

This method blank applies to the following samples: 1045-069-06B, 1045-069-06B DL, 1045-069-06M, 1045-069-06M DL, 1045-069-08M.

DL - Dilution.

J - Indicates an estimated value less than the detection limit.

Client Project ID: LE Carpenter

Job Number: ITPK 51499

### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in Total µg

Sample Matrix: Water

Client Sample ID: Method Blank 3 Lab Sample ID: WB06224

Tentative Identification (1)

None Detected

Concentration (2)

Remarks:

(1) Identification is based on computer search of the NIST Library.

(2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Client Project ID: LE Carpenter

Job Number: ITPK 51499

### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in Total  $\mu g$ Sample Matrix: Water

Client Sample ID: Method Blank 4

Lab Sample ID: WB0623

| Compound                 | Concent | ration     | Compound                  | Concent    | ration   |
|--------------------------|---------|------------|---------------------------|------------|----------|
| acrolein                 | 10      | U          | 1,1-dichloroethene        | 5          | U        |
| acrylonitrile            | 10      | <b>, U</b> | cis-1,2-dichloroethene    | 5          | U        |
| benzene                  | 5       | U          | trans-1,2-dichloroethene  | 5          | U        |
| bromodichloromethane     | 5       | ,<br>U     | 1,2-dichloropropane       | 5          | U        |
| bromoform                | 5       | U          | cis-1,3-dichloropropene   | .5         | U        |
| bromomethane             | 10      | υ          | trans-1,3-dichloropropene | 5          | U        |
| carbon tetrachloride     | 5       | U          | ethyl benzene             | .5         | U        |
| chlorobenzene            | 5       | U          | methylene chloride        | · 1        | J        |
| chloroethane             | 10      | U          | 1,1,2,2-tetrachloroethane | 5          | U        |
| 2-chloroethylvinyl ether | 10      | U          | tetrachloroethene         | 5          | Ü        |
| chloroform               | 5       | U          | toluene                   | 5          | <b>ט</b> |
| chloromethane            | 10      | U          | 1,1,1-trichloroethane     | 5          | U        |
| dibromochloromethane     | 5       | :U         | 1,1,2-trichloroethane     | 5          | U        |
| 1,1-dichloroethane       | 5       | U          | trichloroethene           | · <b>5</b> | U        |
| 1,2-dichloroethane       | 5       | Ŭ          | trichlorofluoromethane    | 5          | U        |
|                          |         | 1 a<br>1 d | vinyl chloride            | 10         | U        |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/23/92

This method blank applies to the following samples: 1045-069-07M DL, 1045-069-08B.

DL - Dilution.

J - Indicates an estimated value less than the detection limit.

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE

KNOXVILLE, TN

Job Number: ITPK 51499

Client Project ID: LE Carpenter

### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in Total µg

Sample Matrix: Water

Client Sample ID: Method Blank 4

Lab Sample ID: WB0623

Tentative Identification (1)

Concentration (2)

None Detected

Remarks:

(1) Identification is based on computer search of the NIST Library.

(2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITPK 51499

# VOST SURROGATE PERCENT RECOVERY SUMMARY

|                  |                   | -         |    | Vo                  | LATILE    |                         |         |
|------------------|-------------------|-----------|----|---------------------|-----------|-------------------------|---------|
| Client Sample ID | Toluen<br>(81-11) |           |    | BF<br><u>(74-12</u> |           | 1,2 Dichloro<br>(70-121 |         |
| 1045-069-05B     | 19                | ajeaje    |    | 29                  | alcole .  | 10                      | aleste. |
| 1045-069-05B DL  | 66                | **        |    | 69                  | **        | 75                      |         |
| 1045-069-05B DL2 | 39                | ***       |    | 39                  | <b>本本</b> | 48                      | **      |
| 1045-069-05M     | 65                | **        |    | 51                  | **        | 72                      | **      |
| 1045-069-05M DL  | 64                | Telephone |    | 54                  | **        | 80                      | ٠       |
| 1045-069-05T     | 26                | **        |    | 18                  | **        | 46                      | alcah   |
| 1045-069-06B     | 86                |           | ٠. | 94                  |           | 95                      |         |
| 1045-069-06B DL  | 91                |           |    | 98                  |           | 115,                    |         |
| 1045-069-06M     | 60                | **        |    | 55                  | **        | 88                      |         |
| 1045-069-06M DL  | 62                | ***       |    | 58                  | <b>in</b> | 96                      |         |
| 1045-069-06T     | 83                | i i       |    | 75                  |           | 88                      |         |
| 1045-069-07M     | 39                | alculu .  |    | 37                  | **        | 83                      |         |
| 1045-069-07M DL  | 47                | ajtajt    |    | 48                  | ***       | 105                     |         |
| 1045-069-07T     | 79.               | **        |    | 80                  |           | · <b>83</b>             |         |
| 1045-069-07T DL  | 57                | **        |    | 68                  | **        | 63                      | *       |
| 1045-069-07T DL2 | 56                | ajeaje    |    | 68                  | **        | 67                      | state   |

DL - Dilution

<sup>\*</sup>Values in parenthesis represent required QC limits.

<sup>\*\*</sup>Values outside of required QC limits.

Job Number: ITPK 51499

# **VOST SURROGATE PERCENT RECOVERY SUMMARY (continued)**

|                  |                          |        | VOLATILE          |                                     |
|------------------|--------------------------|--------|-------------------|-------------------------------------|
| Client Sample ID | Toluene-D8<br>(81-117%)* |        | BFB<br>(74-121%)* | 1,2 Dichloroethane-D4<br>(70-121%)* |
| 1045-069-08B     | 77 **                    |        | 70 **             | 85                                  |
| 1045-069-08M     | 87                       |        | 79                | <b>97</b> .                         |
| 1045-069-08T     | 82                       |        | 71 **             | 87                                  |
| Method Blank 1   | 95                       |        | 89                | 94                                  |
| Method Blank 2   | 108                      | 1      | 97                | 99                                  |
| Method Blank 3   | 97                       | ·<br>• | 96                | 94                                  |
| Method Blank 4   | 101                      |        | 94                | 95                                  |

<sup>\*</sup>Values in parenthesis represent required QC limits.

<sup>\*\*</sup>Values outside of required QC limits.

Job Number: ITPK 51499

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-069-01 Lab Sample ID: K3278

| Compound                   | Concentrat | ion Compound                | Concentration |
|----------------------------|------------|-----------------------------|---------------|
| acenaphthene               | 62 U       | bis(2-chloroisopropyl)ether | 62 U          |
| acenaphthylene             | 62 U       | bis(2-ethylhexyl)phthalate  | 6,500 D       |
| anthracene                 | 62 U       | 4-bromophenyl phenyl ether  | 62 U          |
| benzidine                  | 310 U      | 2-chloronaphthalene         | 62 U          |
| benzo(a)anthracene         | 62 U       | 4-chlorophenyl phenyl ether | 62 U          |
| benzo(b)fluoranthene       | 62 U       | chrysene                    | <b>62</b> U   |
| benzo(k)fluoranthene       | 62 U       | dibenz(a,h)anthracene       | 62 · U        |
| benzo(a)pyrene             | 62 U       | di-n-butylphthalate         | 62 U          |
| benzo(g,h,i)perylene       | 62 U       | 1,2-dichlorobenzene         | 62 U          |
| butylbenzylphthalate       | 62 U       | 1,3-dichlorobenzene         | 62 U          |
| bis(2-chloroethoxy)methane | 62 U       | 1,4-dichlorobenzene         | 62 U          |
| bis(2-chloroethyl)ether    | 62 U       |                             |               |

<sup>-</sup> Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 06/03/92 Date of Analysis:

06/11/92

<sup>-</sup> Indicates an estimated value less than the detection limit.

<sup>-</sup> Compound analyzed at a secondary dilution factor.

Job Number: ITPK 51499

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-069-01 Lab Sample ID: K3278

| Compound                  | Concentration | <u>n</u> | Compound                   | Concentra | <u>ation</u> |
|---------------------------|---------------|----------|----------------------------|-----------|--------------|
| 3,3'dichlorobenzidine     | 120 U         | Ì        | hexachloroethane           | 62        | U            |
| diethylphthalate          | 62 U          |          | indeno(1,2,3-cd)pyrene     | 62        | U            |
| dimethyl phthalate        | 62 U          | ľ.       | isophorone                 | 62        | U            |
| 2,4-dinitrotoluene        | 62 U          |          | naphthalene                | 62        | U            |
| 2,6-dinitrotoluene        | 62 U          | Ī        | nitrobenzene               | 62        | U            |
| di-n-octyl phthalate      | 24 J          |          | n-nitroso-di-n-propylamine | 62        | U            |
| 1,2-diphenylhydrazine(1)  | 62 U          | Ī        | n-nitrosodimethylamine     | 62        | Ü            |
| fluoranthene              | 62 U          | İ        | n-nitrosodiphenylamine(2)  | 62        | U            |
| fluorene                  | 62 U          | Ī        | phenanthrene               | 62        | U            |
| hexachlorobenzene         | 62 U          | i .      | pyrene                     | 62        | U            |
| hexachlorobutadiene       | 62 U          |          | 1,2,4-trichlorobenzene     | 62        | U            |
| hexachlorocyclopentadiene | 62 U          |          |                            |           |              |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

(1) - Screened for as azobenzene

(2) - Detected as diphenylamine

Job Number: ITPK 51499

### ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-069-01 Lab Sample ID: K3278

| Compound                   | Concent | ration       | Compound              | Conce | ntration |
|----------------------------|---------|--------------|-----------------------|-------|----------|
| 4-chloro-3-methylphenol    | 62      | U            | 2-nitrophenol         | 62    | U        |
| 2-chlorophenol             | 62      | <b>ט</b>     | 4-nitrophenol         | 310   | U        |
| 2,4-dichlorophenol         | 62      | U            | pentachlorophenol     | 310   | U        |
| 2,4-dimethylphenol         | 62      | Ü            | phenol                | 62    | U        |
| 2,4-dinitrophenol          | 310     | Ü            | 2,4,6-trichlorophenol | 62    | U        |
| 2-methyl-4,6-dinitrophenol | 310     | <b>U</b> , . |                       |       |          |

Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 06/03/92 Date of Analysis:

06/11/92

Indicates an estimated value less than the detection limit.

Job Number: ITPK 51499

# ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-069-01 Lab Sample ID: K3278

| Tentative Identification (1)          | Concentration (2) |
|---------------------------------------|-------------------|
| 1-dodecanol                           | 8,000             |
| 1-dodecanol                           | 8,300             |
| ethanol, 2-(dodecyloxy)-              | 9 <u>,2</u> 00    |
| 1-hexadecanol                         | 5,400             |
| ethanol, 2-(dodecyloxy)-              | 8,200             |
| ethanol, 2-[(2-ethylhexyl) oxy]-      | 9,300             |
| ethanol, 2-(hexadecyloxy)-            | 5,300             |
| unknown                               | 8,600             |
| 2-undecanol                           | 8,900             |
| unknown                               | 6,400             |
| unknown                               | 8,500             |
| unknown                               | 8,500             |
| unknown                               | 6,400             |
| unknown                               | 8,100             |
| unknown                               | 8,200             |
| unknown                               | 5,400             |
| unknown                               | 7,500             |
| unknown                               | 7,400             |
| unknown                               | 5,200             |
| unknown                               | 5,800             |
| · · · · · · · · · · · · · · · · · · · |                   |

### Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITPK 51499

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-069-02 Lab Sample ID: K3279

| Compound                   | Concent | ration    | Compound                    | Concentr | <u>ation</u> |
|----------------------------|---------|-----------|-----------------------------|----------|--------------|
| acenaphthene               | 10      | U         | bis(2-chloroisopropyl)ether | 10       | U            |
| acenaphthylene             | 10      | <b>U</b>  | bis(2-ethylhexyl)phthalate  | 2        | J            |
| anthracene                 | 10      | U         | 4-bromophenyl phenyl ether  | 10       | U            |
| benzidine                  | 50      | Ü         | 2-chloronaphthalene         | 10       | U            |
| benzo(a)anthracene         | 10      | Ų         | 4-chlorophenyl phenyl ether | 10       | Ü            |
| benzo(b)fluoranthene       | 10      | U         | chrysene                    | 10       | U            |
| benzo(k)fluoranthene       | 10      | U.        | dibenz(a,h)anthracene       | 10       | U            |
| benzo(a)pyrene             | 10      | U         | di-n-butylphthalate         | 10       | U            |
| benzo(g,h,i)perylene       | 10      | U .       | 1,2-dichlorobenzene         | 10       | U            |
| butylbenzylphthalate       | 4       | <b>JB</b> | 1,3-dichlorobenzene         | 10       | U            |
| bis(2-chloroethoxy)methane | 10      | U         | 1,4-dichlorobenzene         | 10       | U            |
| bis(2-chloroethyl)ether    | 10      | U         |                             |          |              |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

KNOXVILLE, TN

Job Number: ITPK 51499

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-069-02 Lab Sample ID: K3279

| Compound                  | Concentration | Compound                   | Concentration |
|---------------------------|---------------|----------------------------|---------------|
| 3,3'dichlorobenzidine     | 20 U          | hexachloroethane           | 10 U          |
| diethylphthalate          | 10 U          | indeno(1,2,3-cd)pyrene     | 10 U          |
| dimethyl phthalate        | 10 U          | isophorone                 | 10 U          |
| 2,4-dinitrotoluene        | 10 U          | naphthalene                | 10 U          |
| 2,6-dinitrotoluene        | 10 Ü          | nitrobenzene               | 10 U          |
| di-n-octyl phthalate      | 10 U          | n-nitroso-di-n-propylamine | 10 U          |
| 1,2-diphenylhydrazine(1)  | 10 U          | n-nitrosodimethylamine     | 10 U          |
| fluoranthene              | 10 U          | n-nitrosodiphenylamine(2)  | 10 U          |
| fluorene                  | 10 U          | phenanthrene               | 10 U          |
| hexachlorobenzene         | 10 U          | pyrene                     | 10 U          |
| hexachlorobutadiene       | 10 U          | 1,2,4-trichlorobenzene     | 10 U          |
| hexachlorocyclopentadiene | 10 U          |                            |               |

- Compound was analyzed for but not detected. The number is the detection limit for the sample.

- Indicates an estimated value less than the detection limit.

(1) - Screened for as azobenzene

(2) - Detected as diphenylamine

Date of Extraction: 06/03/92 Date of Analysis:

06/11/92

Job Number: ITPK 51499

### ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-069-02 Lab Sample ID: K3279

| Compound                   | Concentration | Compound              | Concentration |
|----------------------------|---------------|-----------------------|---------------|
| 4-chloro-3-methylphenol    | 10 U          | 2-nitrophenol         | 10 U          |
| 2-chlorophenol             | 10 U          | 4-nitrophenol         | 50 Ü          |
| 2,4-dichlorophenol         | 10 U          | pentachlorophenol     | 50 U          |
| 2,4-dimethylphenol         | 10 Ü          | phenol                | 10 U          |
| 2,4-dinitrophenol          | 50 U          | 2,4,6-trichlorophenol | 10 U          |
| 2-methyl-4,6-dinitrophenol | 50 U          |                       |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 06/

06/03/92

Date of Analysis:

06/11/92

J - Indicates an estimated value less than the detection limit.

Job Number: ITPK 51499

### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in  $\mu g/\text{liter}$  (ppb)

Sample Matrix: Water

Client Sample ID: 1045-069-02 Lab Sample ID: K3279

| Tentative Identification (1) | Concentration (2) |
|------------------------------|-------------------|
| cyclohexane, isocyanato-     | 14                |
| benzothiazole                | 25                |
| cyclohexane, isothiocyanato- | 76                |
| hydroxylamine, o-decyl-      | 16                |
| unknown                      | 13                |
| trans-2,3-epoxydecane        | 18                |
| unknown (alcohol?)           | 14                |
| unknown                      | 21                |
| unknown                      | 15                |
| unknown                      | 21                |
| unknown                      | 20                |
| unknown                      | 23                |
| unknown (Crown?)             | 120               |
| unknown                      | 42                |
| unknown                      | 220               |
| unknown (Crown?)             | 20                |
| unknown                      | 14                |
| unknown (Crown?)             | 44                |
| unknown                      | 22                |
| unknown                      | 13                |
|                              |                   |

#### Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITPK 51499

### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in  $\mu g$ /liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-069-03 Lab Sample ID: K3280

| Compound                   | Concentration | Compound                    | <b>Concentration</b> |
|----------------------------|---------------|-----------------------------|----------------------|
| acenaphthene               | 10 U          | bis(2-chloroisopropyl)ether | 10 U                 |
| acenaphthylene             | 10 U          | bis(2-ethylhexyl)phthalate  | 10 U                 |
| anthracene                 | 10 U          | 4-bromophenyl phenyl ether  | 10 U                 |
| benzidine                  | 50 U          | 2-chloronaphthalene         | 10 U                 |
| benzo(a)anthracene         | 10 Ŭ          | 4-chlorophenyl phenyl ether | 10 U                 |
| benzo(b)fluoranthene       | 10 U          | chrysene                    | 10 U                 |
| benzo(k)fluoranthene       | 10 U          | dibenz(a,h)anthracene       | 10 Ü                 |
| benzo(a)pyrene             | 10 U -        | di-n-butylphthalate         | 10 U                 |
| benzo(g,h,i)perylene       | 10 U          | 1,2-dichlorobenzene         | 10 U                 |
| butylbenzylphthalate       | 6 JB          | 1,3-dichlorobenzene         | 10 U                 |
| bis(2-chloroethoxy)methane | 10 U          | 1,4-dichlorobenzene         | 10 U                 |
| bis(2-chloroethyl)ether    | 10 U          |                             |                      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

KNOXVILLE, TN

Job Number: ITPK 51499

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-069-03 Lab Sample ID: K3280

| Compound                  | Concentration | Compound                   | <b>Concentration</b> |
|---------------------------|---------------|----------------------------|----------------------|
| 3,3'dichlorobenzidine     | 20 U          | hexachloroethane           | 10 U                 |
| diethylphthalate          | 10 U          | indeno(1,2,3-cd)pyrene     | 10 Ų                 |
| dimethyl phthalate        | 10 U          | isophorone                 | 10 U                 |
| 2,4-dinitrotoluene        | 10 U          | naphthalene                | 10 U                 |
| 2,6-dinitrotoluene        | 10 U          | nitrobenzene               | 10 U                 |
| di-n-octyl phthalate      | 10 U          | n-nitroso-di-n-propylamine | 10 U                 |
| 1,2-diphenylhydrazine(1)  | 10 U          | n-nitrosodimethylamine     | 10 U                 |
| fluoranthene              | 10 U          | n-nitrosodiphenylamine(2)  | 10 · U               |
| fluorene                  | 14 10 U       | phenanthrene               | 10 U                 |
| hexachlorobenzene         | 10 U          | pyrene                     | 10 U                 |
| hexachlorobutadiene       | 10 U          | 1,2,4-trichlorobenzene     | 10 U                 |
| hexachlorocyclopentadiene | 10 U          |                            |                      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

<sup>(1) -</sup> Screened for as azobenzene

<sup>(2) -</sup> Detected as diphenylamine

Job Number: ITPK 51499

# ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in  $\mu g/liter$  (ppb)

Sample Matrix: Water

Client Sample ID: 1045-069-03 Lab Sample ID: K3280

| Compound                   | Concen | tration  | Compound              | Concer | itration |
|----------------------------|--------|----------|-----------------------|--------|----------|
| 4-chloro-3-methylphenol    | 10     | U        | 2-nitrophenol         | 10     | U        |
| 2-chlorophenol             | 10     | Ü        | 4-nitrophenol         | 50     | U        |
| 2,4-dichlorophenol         | 10     | U        | pentachlorophenol     | 50     | U        |
| 2,4-dimethylphenol         | 3      | J        | phenol                | 10     | Ü        |
| 2,4-dinitrophenol          | 50     | <b>ט</b> | 2,4,6-trichlorophenol | 10     | U        |
| 2-methyl-4,6-dinitrophenol | 50     | Ü        |                       |        |          |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Job Number: ITPK 51499

# ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-069-03 Lab Sample ID: K3280

| Tentative Identification (1)          | Concentration (2) |
|---------------------------------------|-------------------|
| 2-pentanone, 4-hydroxy-4-methyl-      | 14 AB             |
| benzothiazole                         | 25                |
| cyclohexane, isothiocyanato-          | 18                |
| 1,1-dodecanediol, diacetate           | 11                |
| decane, 1-fluoro-                     | 13                |
| benzothiazole, 2-(methylthio)-        | 13                |
| trans-2,3-epoxydecane                 | 12                |
| unknown                               | 15                |
| unknown                               | 24                |
| unknown (Crown?)                      | 12                |
| unknown                               | 26                |
| 1,4,7,10,13,16-hexaoxacyclooctadecane | 12                |
| unknown                               | 13                |
| unknown (Crown?)                      | 49                |
| 1,4,7,10,13,16-hexaoxacyclooctadecane | 17                |
| unknown                               | 16                |
| unknown                               | 14                |
| unknown                               | 42                |
| 1,4,7,10,13,16-hexaoxacyclooctadecane | 14                |
| 1,4,7,10,13,16-hexaoxacyclooctadecane | 26                |

#### Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- A Suspected aldol condensation product.
- B Analyte was found in the blank as well as the sample.

Job Number: ITPK 51499

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-069-04 Lab Sample ID: K3281

| Compound                   | Concen | tration  | Compound                    | Concentr | ation    |
|----------------------------|--------|----------|-----------------------------|----------|----------|
| acenaphthene               | 50     | U        | bis(2-chloroisopropyl)ether | 50       | U        |
| acenaphthylene             | 50     | U        | bis(2-ethylhexyl)phthalate  | 38       | J        |
| anthracene                 | 50     | U        | 4-bromophenyl phenyl ether  | 50       | U        |
| benzidine                  | 250    | U        | 2-chloronaphthalene         | 50       | <b>บ</b> |
| benzo(a)anthracene         | 50     | Ü.       | 4-chlorophenyl phenyl ether | 50       | Ü        |
| benzo(b)fluoranthene       | 50     | U.       | chrysene                    | 50       | U        |
| benzo(k)fluoranthene       | 50     | U        | dibenz(a,h)anthracene       | 50       | U        |
| benzo(a)pyrene             | 50     | U        | di-n-butylphthalate         | 50       | U        |
| benzo(g,h,i)perylene       | 50     | U        | 1,2-dichlorobenzene         | 50       | Ü        |
| butylbenzylphthalate       | 50     | U        | 1,3-dichlorobenzene         | 50       | U        |
| bis(2-chloroethoxy)methane | 50     | <b>U</b> | 1,4-dichlorobenzene         | 50       | U        |
| bis(2-chloroethyl)ether    | 50     | U.       |                             |          |          |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Job Number: ITPK 51499

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-069-04 Lab Sample ID: K3281

| Compound                  | <u>Concentration</u> | Compound                   | Concentration |
|---------------------------|----------------------|----------------------------|---------------|
| 3,3'dichlorobenzidine     | 100 U                | hexachloroethane           | <b>50</b> U   |
| diethylphthalate          | .50 U                | indeno(1,2,3-cd)pyrene     | . 50 U        |
| dimethyl phthalate        | <i>5</i> 0 Ū         | isophorone                 | <b>50</b> U   |
| 2,4-dinitrotoluene        | 50 U                 | naphthalene                | <b>50</b> U   |
| 2,6-dinitrotoluene        | <b>50 U</b>          | nitrobenzene               | 50 Ü          |
| di-n-octyl phthalate      | 50 U                 | n-nitroso-di-n-propylamine | 50 U          |
| 1,2-diphenylhydrazine(1)  | <b>50</b> U          | n-nitrosodimethylamine     | 50 U          |
| fluoranthene              | 50 U                 | n-nitrosodiphenylamine(2)  | 50 U          |
| fluorene                  | 50 Ü                 | phenanthrene               | <b>50</b> U   |
| hexachlorobenzene         | <b>50</b> Ü          | pyrene                     | 50 U          |
| hexachlorobutadiene       | 50 U                 | 1,2,4-trichlorobenzene     | 50 U          |
| hexachlorocyclopentadiene | <b>50</b> U          |                            |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

(1) - Screened for as azobenzene(2) - Detected as diphenylamine

Client Project ID: LE Carpenter

Job Number: ITPK 51499

# ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in  $\mu$ g/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-069-04 Lab Sample ID: K3281

| Compound                   | Concer | tration | Compound              | Conce | ntration |
|----------------------------|--------|---------|-----------------------|-------|----------|
| 4-chloro-3-methylphenol    | 50     | U       | 2-nitrophenol         | 50    | U        |
| 2-chlorophenol             | 50     | U       | 4-nitrophenol         | 250   | U.       |
| 2,4-dichlorophenol         | 50     | U       | pentachlorophenol     | 250   | U        |
| 2,4-dimethylphenol         | 50     | U :     | phenol                | 50    | U        |
| 2,4-dinitrophenol          | 250    | U       | 2,4,6-trichlorophenol | 50    | U        |
| 2-methyl-4,6-dinitrophenol | 250    | U       |                       |       |          |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

Job Number: ITPK 51499

### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-069-04 Lab Sample ID: K3281

| Tentative Identification (1)                         | Concentration (2) |
|--|-------------------|
| hexanal  | 190               |
| heptanal   | 290               |
| 2-nonenal, (E)-                                      | 210               |
| octanoic acid  | 260               |
| cyclopropane, 1-heptyl-2-methyl-                     | 260               |
| nonanoic acid  | 210               |
| 1-dodecanol  | 290               |
| decanoic acid  | 220               |
| dodecanoic acid                                      | 680               |
| tetradecanoic acid                                   | 970               |
| 9-hexadecenoic acid                                  | 970               |
| hexadecanoic acid                                    | 920               |
| 2-octadecenal  | 4,000             |
| 9-octadecenoic acid (Z)-, 2,-3-dihydroxypropyl ester | 720               |
| unknown  | 1,100             |
| 1,4,7,10,13,16-hexaoxacyclooctadecane                | 230               |
| unknown  | 890               |
| unknown (Ester?)                                     | 220               |
| unknown (Ester?)                                     | 1,300             |
| unknown  | 450               |

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITPK 51499

#### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-069-04 RE Lab Sample ID: K3281 RE

| Compound                   | Concentration | Compound                    | Concentration |
|----------------------------|---------------|-----------------------------|---------------|
| acenaphthene               | 266 U         | bis(2-chloroisopropyl)ether | 266 U         |
| acenaphthylene             | 266 U         | bis(2-ethylhexyl)phthalate  | 90 J          |
| anthracene                 | 266 U         | 4-bromophenyl phenyl ether  | 266 U         |
| benzidine                  | 665 U         | 2-chloronaphthalene         | 266 U         |
| benzo(a)anthracene         | 266 U         | 4-chlorophenyl phenyl ether | 266 U         |
| benzo(b)fluoranthene       | 266 U         | chrysene                    | 266 U         |
| benzo(k)fluoranthene       | 266 U         | dibenz(a,h)anthracene       | 266 U         |
| benzo(a)pyrene             | 266 U         | di-n-butylphthalate         | 266 U         |
| benzo(g,h,i)perylene       | 266 U         | 1,2-dichlorobenzene         | 266 U         |
| butylbenzylphthalate       | 266 U         | 1,3-dichlorobenzene         | 266 Ü         |
| bis(2-chloroethoxy)methane | 266 U         | 1,4-dichlorobenzene         | 266 U         |
| bis(2-chloroethyl)ether    | 266 U         |                             |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 06/24/92 Date of Analysis: 06/26/92

J - Indicates an estimated value less than the detection limit.

Job Number: ITPK 51499

# BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-069-04 RE Lab Sample ID: K3281 RE

| Compound                  | Concentration | Compound                   | Concentration |
|---------------------------|---------------|----------------------------|---------------|
| 3,3'dichlorobenzidine     | 665 U         | hexachloroethane           | 266 U         |
| diethylphthalate          | 266 U         | indeno(1,2,3-cd)pyrene     | 266 U         |
| dimethyl phthalate        | 266 U         | isophorone                 | 266 U         |
| 2,4-dinitrotoluene        | 266 U         | naphthalene                | 266 U         |
| 2,6-dinitrotoluene        | 266 U         | nitrobenzene               | 266 U         |
| di-n-octyl phthalate      | 266 U         | n-nitroso-di-n-propylamine | 266 U         |
| 1,2-diphenylhydrazine(1)  | 266 U         | n-nitrosodimethylamine     | 266 U         |
| fluoranthene              | 266 U         | n-nitrosodiphenylamine(2)  | 266 U         |
| fluorene                  | 266 U         | phenanthrene               | 266 Ü         |
| hexachlorobenzene         | 266 U         | pyrene                     | 266 U         |
| hexachlorobutadiene       | 266 U         | 1,2,4-trichlorobenzene     | 266 U         |
| hexachlorocyclopentadiene | 266 Ū         |                            |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

(1) - Screened for as azobenzene

(2) - Detected as diphenylamine

Date of Extraction:

06/24/92

Date of Analysis:

06/26/92

Job Number: ITPK 51499

#### ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-069-04 RE Lab Sample ID: K3281 RE

| Compound                   | Concen | tration  | Compound              | Conce | ntration |
|----------------------------|--------|----------|-----------------------|-------|----------|
| 4-chloro-3-methylphenol    | 266    | U .      | 2-nitrophenol         | 266   | U ·      |
| 2-chlorophenol             | 266    | U        | 4-nitrophenol         | 665   | U        |
| 2,4-dichlorophenol         | 266    | U        | pentachlorophenol     | 665   | U        |
| 2,4-dimethylphenol         | 266    | U .      | phenol                | 266   | U        |
| 2,4-dinitrophenol          | 665    | Ų        | 2,4,6-trichlorophenol | 266   | U        |
| 2-methyl-4,6-dinitrophenol | 665    | <b>u</b> |                       |       |          |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 06/24/92 Date of Analysis: 06/26/92

J - Indicates an estimated value less than the detection limit.

Job Number: ITPK 51499

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: 1045-069-04 RE Lab Sample ID: K3281 RE

| Tentative Identification (1)          | Concentration (2) |
|---------------------------------------|-------------------|
| dodecanoic acid                       | 3,700             |
| tetradecanoic acid                    | 8,400             |
| 9-hexadecenoic acid                   | 12,000            |
| hexadecanoic acid                     | 12,000            |
| unknown                               | 3,500             |
| ethanol, 2,2'-[oxybis(2,1-ET          | 2,800             |
| unknown                               | 38,000            |
| 15-crown-5                            | 10,000            |
| unknown                               | 3,100             |
| unknown                               | 5,500             |
| 15-crown-5                            | 17,000            |
| unknown                               | 4,000             |
| unknown                               | 3,000             |
| 1,4,7,10,13,16-hexaoxacyclooctadecane | 26,000            |
| unknown                               | 6,100             |
| unknown                               | 6,400             |
| 1,4,7,10,13,16-hexaoxacyclooctadecane | 16,000            |
| unknown                               | 4,500             |
| unknown                               | 4,100             |
| 1,4,7,10,13,16-hexaoxacyclooctadecane | 4,600             |
|                                       |                   |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

KNOXVILLE, TN

Job Number: ITPK 51499

#### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 1

Lab Sample ID: BL0863

| Compound                   | Concentration | Compound                    | Concentration |
|----------------------------|---------------|-----------------------------|---------------|
| acenaphthene               | 10 U          | bis(2-chloroisopropyl)ether | 10 U          |
| acenaphthylene             | 10 U          | bis(2-ethylhexyl)phthalate  | 10 U          |
| anthracene                 | 10 U          | 4-bromophenyl phenyl ether  | 10 U          |
| benzidine                  | .50 Ū         | 2-chloronaphthalene         | 10 U          |
| benzo(a)anthracene         | 10 U          | 4-chlorophenyl phenyl ether | Ì0 U          |
| benzo(b)fluoranthene       | 10 U          | chrysene                    | 10 Ü          |
| benzo(k)fluoranthene       | 10 U          | dibenz(a,h)anthracene       | 10 U          |
| benzo(a)pyrene             | 10 U          | di-n-butylphthalate         | 10 U          |
| benzo(g,h,i)perylene       | 10 U          | 1,2-dichlorobenzene         | 10 Ü          |
| butylbenzylphthalate       | <b>3 J</b>    | 1,3-dichlorobenzene         | 10 Ü          |
| bis(2-chloroethoxy)methane | 10 U          | 1,4-dichlorobenzene         | 10 U          |
| bis(2-chloroethyl)ether    | .10 U         | t -                         |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 06/03/92 Date of Analysis: 06/11/92

J - Indicates an estimated value less than the detection limit.

Job Number: ITPK 51499

### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in  $\mu g$ /liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 1

Lab Sample ID: BL0863

| Compound                  | Concentra | ation . | Compound                   | Concentra | <u>ation</u> |
|---------------------------|-----------|---------|----------------------------|-----------|--------------|
| 3,3'dichlorobenzidine     | 20        | U       | hexachloroethane           | 10        | U            |
| diethylphthalate          | 10        | U       | indeno(1,2,3-cd)pyrene     | 10        | U            |
| dimethyl phthalate        | 10        | Ü       | isophorone                 | . 10      | U            |
| 2,4-dinitrotoluene        | 10        | Ü       | naphthalene                | 10        | U            |
| 2,6-dinitrotoluene        | 10        | U       | nitrobenzene               | 10        | U            |
| di-n-octyl phthalate      | 10        | U       | n-nitroso-di-n-propylamine | 10        | U            |
| 1,2-diphenylhydrazine(1)  | 10        | U       | n-nitrosodimethylamine     | 10        | <b>U</b> .   |
| fluoranthene              | 10        | U       | n-nitrosodiphenylamine(2)  | 10        | Ū            |
| fluorene                  | 10        | U       | phenanthrene               | 10        | Ü            |
| hexachlorobenzene         | . 10      | U       | pyrene                     | 10        | U            |
| hexachlorobutadiene       | 10        | U       | 1,2,4-trichlorobenzene     | 10        | Ù            |
| hexachlorocyclopentadiene | 10        | U       |                            |           |              |

- U Compound was analyzed for but not detected. The number is the detection limit for the sample.
- J Indicates an estimated value less than the detection limit.
- (1) Screened for as azobenzene
- (2) Detected as diphenylamine

Date of Extraction:

Date of Analysis:

06/03/92

06/11/92

This method blank applies to the following samples: 1045-069-01, 1045-069-02, 1045-069-03 and 1045-069-04.

Job Number: ITPK 51499

#### ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 1

Lab Sample ID: BL0863

| Compound                   | Concentration | Compound              | Concentration |
|----------------------------|---------------|-----------------------|---------------|
| 4-chloro-3-methylphenol    | 10 U          | 2-nitrophenol         | 10 U          |
| 2-chlorophenol             | 10 U          | 4-nitrophenol         | 50 U          |
| 2,4-dichlorophenol         | 10 U          | pentachlorophenol     | <b>50</b> U   |
| 2,4-dimethylphenol         | 10 U          | phenol                | 10 U          |
| 2,4-dinitrophenol          | 50 U          | 2,4,6-trichlorophenol | 10 U          |
| 2-methyl-4,6-dinitrophenol | <b>50</b> U   |                       | •             |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 06/03/92 Date of Analysis: 06/11/92

J - Indicates an estimated value less than the detection limit.

Client Project ID: LE Carpenter

Job Number: ITPK 51499

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 1

Lab Sample ID: BL0863

Tentative Identification (1)

Concentration (2)

2-pentanone, 4-hydroxy-4-methyl-

36 A

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- A Suspected aldol condensation product.

Job Number: ITPK 51499

#### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in  $\mu$ g/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 2

Lab Sample ID: BL0840

| Compound                   | Concentration | Compound                    | Concentration |
|----------------------------|---------------|-----------------------------|---------------|
| acenaphthene               | 10 U          | bis(2-chloroisopropyl)ether | 10 U          |
| acenaphthylene             | 10 U          | bis(2-ethylhexyl)phthalate  | 10 U          |
| anthracene                 | 10 U          | 4-bromophenyl phenyl ether  | 10 U          |
| benzidine                  | 25 U          | 2-chloronaphthalene         | 10 U          |
| benzo(a)anthracene         | 10 U          | 4-chlorophenyl phenyl ether | 10 U          |
| benzo(b)fluoranthene       | 10 Ŭ          | chrysene                    | 10 U          |
| benzo(k)fluoranthene       | 10 U          | dibenz(a,h)anthracene       | 10 U          |
| benzo(a)pyrene             | 10 U          | di-n-butylphthalate         | 10 U          |
| benzo(g,h,i)perylene       | 10 U          | 1,2-dichlorobenzene         | 10 U          |
| butylbenzylphthalate       | 1 J           | 1,3-dichlorobenzene         | 10 U          |
| bis(2-chloroethoxy)methane | 10 Ù          | 1,4-dichlorobenzene         | 10 U          |
| bis(2-chloroethyl)ether    | 10 U          |                             |               |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 06/24/92 Date of Analysis: 06/26/92

J - Indicates an estimated value less than the detection limit.

Job Number: ITPK 51499

### BASE/NEUTRAL EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS (continued)

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 2

Lab Sample ID: BL0840

| Compound                  | Concentration | Compound                   | <u>Concentration</u> |
|---------------------------|---------------|----------------------------|----------------------|
| 3,3'dichlorobenzidine     | 25 U          | hexachloroethane           | 10 <u>U</u>          |
| diethylphthalate          | 10 U          | indeno(1,2,3-cd)pyrene     | . 10 U               |
| dimethyl phthalate        | 10 U          | isophorone                 | 10 U                 |
| 2,4-dinitrotoluene        | 10 U          | naphthalene                | . 10 Ü               |
| 2,6-dinitrotoluene        | 10 U          | nitrobenzene               | 10 U                 |
| di-n-octyl phthalate      | 10 U          | n-nitroso-di-n-propylamine | 10 U                 |
| 1,2-diphenylhydrazine(1)  | 10 U          | n-nitrosodimethylamine     | 10 U                 |
| fluoranthene              | 10 U          | n-nitrosodiphenylamine(2)  | 10 U                 |
| fluorene                  | 10 U          | phenanthrene               | 10 U                 |
| hexachlorobenzene         | 10 U          | pyrene                     | 10 U                 |
| hexachlorobutadiene       | 10 U          | 1,2,4-trichlorobenzene     | 10 U                 |
| hexachlorocyclopentadiene | 10 U          |                            |                      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

J - Indicates an estimated value less than the detection limit.

(1) - Screened for as azobenzene

(2) - Detected as diphenylamine

Date of Extraction: 06/24/92 Date of Analysis: 06/26/92

This method blank applies to the following samples: 1045-069-04 RE.

Job Number: ITPK 51499

#### ACID EXTRACTABLE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 2

Lab Sample ID: BL0840

| Compound                   | Concent | ration | Compound              | Concer | ntration |
|----------------------------|---------|--------|-----------------------|--------|----------|
| 4-chloro-3-methylphenol    | 10      | U      | 2-nitrophenol         | .10    | U        |
| 2-chlorophenol             | 10      | U      | 4-nitrophenol         | 25     | Ū        |
| 2,4-dichlorophenol         | 10      | U      | pentachlorophenol     | 25     | U        |
| 2,4-dimethylphenol         | 10      | U      | phenol                | 10     | U        |
| 2,4-dinitrophenol          | 25      | U · ;  | 2,4,6-trichlorophenol | 10     | U .      |
| 2-methyl-4,6-dinitrophenol | 25      | U .    |                       |        |          |

<sup>-</sup> Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Extraction: 06/24/92 Date of Analysis:

06/26/92

<sup>-</sup> Indicates an estimated value less than the detection limit.

Client Project ID: LE Carpenter

Job Number: ITPK 51499

#### ADDITIONAL SEMIVOLATILE ORGANIC COMPOUNDS

Results in µg/liter (ppb)

Sample Matrix: Water

Client Sample ID: Method Blank 2

Lab Sample ID: BL0840

| Tentative Identification (1)  | Concentration (2) |
|-------------------------------|-------------------|
| unknown                       | 28                |
| benzene, 1, 1'-sulfonylbis[4- | 26                |
| acetic acid, (triphenylphosp  | 68                |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITPK 51499

#### WATER SURROGATE PERCENT RECOVERY SUMMARY

#### **SEMI-VOLATILE**

| Client Sample ID | Nitro-<br>Benzene-D5<br>(35-114%)* | 2-Fluoro-<br>Biphenyl<br>(43-116%)* | Terphenyl-<br>D14<br>(33-141%)* | Phenol-D5<br>(10-94%)* | 2-Fluoro-<br>Phenol<br>(21-100%)* | 2,4,6-<br>Tribromo-<br>Phenol<br>(10-123%)* |
|------------------|------------------------------------|-------------------------------------|---------------------------------|------------------------|-----------------------------------|---|
| 1045-069-01      | 57                                 | 54                                  | 49                              | 42                     | 50                                | 32  |
| 1045-069-01 DL   | D                                  | D                                   | D                               | D                      | Ď                                 | D   |
| 1045-069-02      | 62                                 | 69                                  | 81                              | 33                     | 47                                | 77  |
| 1045-069-03      | 76                                 | 58                                  | 87                              | 34                     | 50                                | 75  |
| 1045-069-04      | 30 ***                             | 34 **                               | 38                              | 31                     | 46                                | 73  |
| Method Blank 1   | 77                                 | 85                                  | 94                              | 33                     | 53                                | 75  |

<sup>\*</sup>Values in parenthesis represent required QC limits.

<sup>\*\*</sup>Values outside required QC limits.

D - Surrogates diluted out.

DL - Dilution.

Job Number: ITPK 51499

Client Project ID: LE Carpenter

#### WATER SURROGATE PERCENT RECOVERY SUMMARY

#### **SEMI-VOLATILE**

| Client Sample ID | Nitro-<br>Benzene-D5<br>(35-114%)* | 2-Fluoro-<br>Biphenyl<br>(43-116%)* | Terphenyl-<br>D14<br>(33-141%)* | Phenol-D5<br>(10-94%)* | 2-Fluoro-<br>Phenol<br>(21-100%)* | 2,4,6-<br>Tribromo-<br>Phenol<br>(10-123%)* |
|------------------|------------------------------------|-------------------------------------|---------------------------------|------------------------|-----------------------------------|---|
| 1045-069-04 RE   | 79                                 | 94                                  | 92                              | 102 **                 | 96                                | 88  |
| Method Blank 2   | 78                                 | 83                                  | 86                              | <b>77</b> .            | 74                                | 85  |

<sup>\*</sup>Values in parenthesis represent required QC limits.

<sup>\*\*</sup>Values outside of required QC limits.

### Appendix G

Tenax Resin Analyses

Client Project ID: LE Carpenter Job Number: ITPK 51499

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in Total  $\mu g$ Sample Matrix: VOST Tube

Client Sample ID: 1045-069-05T Lab Sample ID: K3270T

| Compound                 | Concentration       | Compound                  | Concentration |  |
|--------------------------|---------------------|---------------------------|---------------|--|
| acrolein                 | 10 U                | 1,1-dichloroethene        | 5 Ü           |  |
| acrylonitrile            | 10 U                | cis-1,2-dichloroethene    | 5 U           |  |
| benzene                  | 5 · U               | trans-1,2-dichloroethene  | 5 U           |  |
| bromodichloromethane     | 5 Ü                 | 1,2-dichloropropane       | 5 U           |  |
| bromoform                | 5 U                 | cis-1,3-dichloropropene   | 5 U           |  |
| bromomethane             | 10 U                | trans-1,3-dichloropropene | 5 U           |  |
| carbon tetrachloride     | 5 U                 | ethyl benzene             | , 5 U         |  |
| chlorobenzene            | 5 U                 | methylene chloride        | 8 B           |  |
| chloroethane             | 10 U                | 1,1,2,2-tetrachloroethane | 5 U           |  |
| 2-chloroethylvinyl ether | 10 U                | tetrachloroethene         | 5 U           |  |
| chloroform               | <b>5 U</b>          | toluene                   | 5 U           |  |
| chloromethane            | 10 U                | 1,1,1-trichloroethane     | 5 U           |  |
| dibromochloromethane     | <b>5</b> , <b>U</b> | 1,1,2-trichloroethane     | 5 Ü           |  |
| 1,1-dichloroethane       | 5 U                 | trichloroethene           | 5 U           |  |
| 1,2-dichloroethane       | 5 U                 | trichlorofluoromethane    | 5 U           |  |
|                          |                     | vinyl chloride            | 10 U          |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/19/92

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

Job Number: ITPK 51499

Client Project ID: LE Carpenter

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in Total µg

Sample Matrix: VOST Tube

Client Sample ID: 1045-069-05T Lab Sample ID: K3270T

Tentative Identification (1)

Concentration (2)

Unknown

5.3

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Client Project ID: LE Carpenter

Job Number: ITPK 51499

### VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS

Results in Total  $\mu g$ Sample Matrix: VOST Tube

Client Sample ID: 1045-069-05M Lab Sample ID: K3270MR

| Compound                 | Concen | tration | Compound                  | Concentr |    |
|--------------------------|--------|---------|---------------------------|----------|----|
| acrolein                 | 10     | Ŭ       | 1,1-dichloroethene        | 5        | U  |
| acrylonitrile            | 10     | U       | cis-1,2-dichloroethene    | .5       | U  |
| benzene                  | 6      | ;       | trans-1,2-dichloroethene  | 5        | Ü  |
| bromodichloromethane     | 5      | U       | 1,2-dichloropropane       | 5        | U  |
| bromoform                | 5      | U       | cis-1,3-dichloropropene   | 5        | Ų  |
| bromomethane             | 10     | U       | trans-1,3-dichloropropene | 5        | U  |
| carbon tetrachloride     | 5      | υ       | ethyl benzene             | 5        | Ü  |
| chlorobenzene            | 5      | Ü       | methylene chloride        | 380      | BD |
| chloroethane             | 10     | U       | 1,1,2,2-tetrachloroethane | 5        | U  |
| 2-chloroethylvinyl ether | 10     | U       | tetrachloroethene         | 5        | U  |
| chloroform               | 7      |         | toluene                   | 5        | U  |
| chloromethane            | 1.0    | U       | 1,1,1-trichloroethane     | 5        | U  |
| dibromochloromethane     | 5      | U       | 1,1,2-trichloroethane     | 5        | U  |
| 1,1-dichloroethane       | .5     | U       | trichloroethene           | . 5      | Ü  |
| 1,2-dichloroethane       | 5      | U       | trichlorofluoromethane    | 5        | U  |
|                          |        |         | vinyl chloride            | 10       | ·U |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/22/92

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

D - Compound analyzed at a secondary dilution factor.

Client Project ID: LE Carpenter

Job Number: ITPK 51499

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in Total µg

Sample Matrix: VOST Tube

Client Sample ID: 1045-069-05M Lab Sample ID: K3270MR

| Concentration (2) |
|-------------------|
| 15                |
| 5.4               |
| 6.6               |
|                   |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Client Project ID: LE Carpenter

Job Number: ITPK 51499

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in Total μg
Sample Matrix: VOST Tube

Client Sample ID: 1045-069-05B Lab Sample ID: K3270B

| Compound                 | Concentration | Compound                  | Concentration |            |
|--------------------------|---------------|---------------------------|---------------|------------|
| acrolein                 | 10 Ü          | 1,1-dichloroethene        | 5             | U          |
| acrylonitrile            | 10 U          | cis-1,2-dichloroethene    | 5             | · <b>U</b> |
| benzene                  | 48            | trans-1,2-dichloroethene  | 5             | U          |
| bromodichloromethane     | 5 U           | 1,2-dichloropropane       | 5             | Ü          |
| bromoform                | 5 U           | cis-1,3-dichloropropene   | . 5           | U          |
| bromomethane             | 10 U          | trans-1,3-dichloropropene | . 5           | U          |
| carbon tetrachloride     | 5 U           | ethyl benzene             | 430           | Ď          |
| chlorobenzene            | 5 U           | methylene chloride        | 170           | В          |
| chloroethane             | 10 s. U       | 1,1,2,2-tetrachloroethane | 5             | U          |
| 2-chloroethylvinyl ether | 10 U          | tetrachloroethene         | 5             | U          |
| chloroform               | 15            | toluene                   | 140           | D          |
| chloromethane            | 10 U          | 1,1,1-trichloroethane     | 5             | U          |
| dibromochloromethane     | <b>5</b> U    | 1,1,2-trichloroethane     | 5             | U          |
| 1,1-dichloroethane       | 5 U           | trichloroethene           | 5             | U          |
| 1,2-dichloroethane       | <b>5</b> U    | trichlorofluoromethane    | 5             | · <b>U</b> |
|                          |               | vinyl chloride            | 10            | U          |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/19/92

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

D - Compound analyzed at a secondary dilution factor.

Job Number: ITPK 51499

Client Project ID: LE Carpenter

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in Total µg

Sample Matrix: VOST Tube

Client Sample ID: 1045-069-05B Lab Sample ID: K3270B

| Tentative Identification (1)    | Concentration (2) |  |  |
|---------------------------------|-------------------|--|--|
| benzene, dimethyl-              | 240 Y             |  |  |
| cyclopentane, 1-hexyl-3-methyl- | 21                |  |  |
| undecane, 5,6-dimethyl-         | 29                |  |  |
| octane, 3,5-dimethyl-           | 61                |  |  |
| nonane, 3-methyl-               | 37                |  |  |
| benzene, ethylmethyl-           | 61 Y              |  |  |
| 2-dodecen-1-oL, 12-chloro-      | 40                |  |  |
| decane, 2,5,6-trimethyl-        | 44                |  |  |
| benzene, trimethyl-             | 47 Y              |  |  |
| undecane                        | 43                |  |  |

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Y - Indistinguishable isomer in tentatively identified compounds.

Job Number: ITPK 51499

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in Total  $\mu g$ Sample Matrix: VOST Tube

Client Sample ID: 1045-069-06T Lab Sample ID: K3271T

| Compound                 | Concentration | Compound                  | Concentration |           |
|--------------------------|---------------|---------------------------|---------------|-----------|
| acrolein                 | 10 U          | 1,1-dichloroethene        | 5             | U         |
| acrylonitrile            | 10 U          | cis-1,2-dichloroethene    | 5             | U         |
| benzene                  | 5 U           | trans-1,2-dichloroethene  | .5            | U         |
| bromodichloromethane     | 5 U           | 1,2-dichloropropane       | 5             | Ų         |
| bromoform                | 5 U           | cis-1,3-dichloropropene   | . 5           | U         |
| bromomethane             | 10 · U        | trans-1,3-dichloropropene | 5             | U         |
| carbon tetrachloride     | 5 Ü           | ethyl benzene             | 5             | U         |
| chlorobenzene            | .5 U          | methylene chloride        | 3             | JB        |
| chloroethane             | 10 U          | 1,1,2,2-tetrachloroethane | 5             | U         |
| 2-chloroethylvinyl ether | 10 U          | tetrachloroethene         | 5             | Ü         |
| chloroform               | 5 U           | toluene                   | 1             | <b>.J</b> |
| chloromethane            | 10 U          | 1,1,1-trichloroethane     | 5             | U         |
| dibromochloromethane     | 5 U           | 1,1,2-trichloroethane     | 5             | U         |
| 1,1-dichloroethane       | 5 U           | trichloroethene           | 5             | ับ        |
| 1,2-dichloroethane       | 5 U           | trichlorofluoromethane    | 5             | U         |
|                          |               | vinyl chloride            | 10            | Ų         |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/19/92

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

June 30, 1992

Client Project ID: LE Carpenter

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

Job Number: ITPK 51499

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in Total µg

Sample Matrix: VOST Tube

Client Sample ID: 1045-069-06T Lab Sample ID: K3271T

Tentative Identification (1)

Concentration (2)

None Detected

Remarks:

(1) Identification is based on computer search of the NIST Library.

(2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Client Project ID: LE Carpenter

Job Number: ITPK 51499

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in Total µg
Sample Matrix: VOST Tube

Client Sample ID: 1045-069-06M Lab Sample ID: K3271M

| Compound                 | Concentration           | on (           | Compound                  | Concent    | ration |
|--------------------------|-------------------------|----------------|---------------------------|------------|--------|
| acrolein                 | 10 U                    | 1              | ,1-dichloroethene         | 5          | U      |
| acrylonitrile            | 10 U                    | c              | is-1,2-dichloroethene     | 5          | U      |
| benzene                  | <b>5</b> ' ' <b>U</b> ' | t              | rans-1,2-dichloroethene   | 5          | U      |
| bromodichloromethane     | 5 U                     | 1              | ,2-dichloropropane        | 5          | U      |
| bromoform                | <b>5</b> :: <b>U</b>    | · , c          | sis-1,3-dichloropropene   | .5         | U      |
| bromomethane             | 10 U                    | t              | rans-1,3-dichloropropene  | 5          | U      |
| carbon tetrachloride     | 5 U                     | •              | ethyl benzene             | 5          | U      |
| chlorobenzene            | 5 U                     | ·. r           | nethylene chloride        | 240        | DB     |
| chloroethane             | 10 U                    | 1              | 1,1,2,2-tetrachloroethane | 5          | U      |
| 2-chloroethylvinyl ether | 10 U                    | , , <b>, t</b> | etrachloroethene          | 5          | U      |
| chloroform               | <b>2</b> J              | t              | oluene                    | 5          | U .    |
| chloromethane            | 10 U                    | 1              | 1,1,1-trichloroethane     | <b>.</b> 5 | U      |
| dibromochloromethane     | 5 U                     | 1              | 1,1,2-trichloroethane     | 5          | U      |
| 1,1-dichloroethane       | 5 U                     | t              | richloroethene            | 5          | U      |
| 1,2-dichloroethane       | 5 Ü                     | .t             | richlorofluoromethane     | 5          | U      |
|                          |                         | 10.0           | rinyl chloride            | 10         | Ū      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/22/92

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

D - Compound analyzed at a secondary dilution factor.

Client Project ID: LE Carpenter

Job Number: ITPK 51499

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in Total µg

Sample Matrix: VOST Tube

Client Sample ID: 1045-069-06M Lab Sample ID: K3271M

Tentative Identification (1)

Concentration (2)

None Detected

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITPK 51499

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in Total  $\mu g$ Sample Matrix: VOST Tube

Client Sample ID: 1045-069-06B Lab Sample ID: K3271B

| Compound                 | Concentration |    | Compound                  | Concent    | Concentration |  |
|--------------------------|---------------|----|---------------------------|------------|---------------|--|
| acrolein                 | 10            | U  | 1,1-dichloroethene        | 5          | U             |  |
| acrylonitrile            | 10            | Ü  | cis-1,2-dichloroethene    | 5          | U             |  |
| benzene                  | 130           |    | trans-1,2-dichloroethene  | 5          | Ü             |  |
| bromodichloromethane     | 2             | J  | 1,2-dichloropropane       | 5          | U             |  |
| bromoform                | <b>5</b>      | U  | cis-1,3-dichloropropene   | 5          | U             |  |
| bromomethane             | 10            | Ü  | trans-1,3-dichloropropene | 5          | U             |  |
| carbon tetrachloride     | 5             | U  | ethyl benzene             | 62         |               |  |
| chlorobenzene            | 5             | U  | methylene chloride        | 450        | DB            |  |
| chloroethane             | 10.1          | U. | 1,1,2,2-tetrachloroethane | 5          | Ü             |  |
| 2-chloroethylvinyl ether | 10            | U  | tetrachloroethene         | 5          | U             |  |
| chloroform               | 30            |    | toluene                   | 360        | D             |  |
| chloromethane            | 10            | U  | 1,1,1-trichloroethane     | 2          | J             |  |
| dibromochloromethane     | 5             | U  | 1,1,2-trichloroethane     | 5          | · <b>U</b>    |  |
| 1,1-dichloroethane       | 5             | U  | trichloroethene           | · <b>5</b> | Ų             |  |
| 1,2-dichloroethane       | 5             | Ų  | trichlorofluoromethane    | 5          | U             |  |
|                          |               |    | vinyl chloride            | 10         | U             |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/22/92

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

D - Compound analyzed at a secondary dilution factor.

Job Number: ITPK 51499

Client Project ID: LE Carpenter

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in Total µg

Sample Matrix: VOST Tube

Client Sample ID: 1045-069-06B Lab Sample ID: K3271B

| Tentative Identification (1) | Concentration (2) |
|------------------------------|-------------------|
| ethane, 1,1,2-trichloro-1,2, | 8.6               |
| benzene, dimethyl-           | 1,300 Y           |
| benzene, 1,2-dimethyl-       | 180               |

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Y - Indistinguishable isomer in tentatively identified compounds.

Job Number: ITPK 51499

### VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS Results in Total µg

Sample Matrix: VOST Tube

Client Sample ID: 1045-069-07T Lab Sample ID: K3272T

| Compound                 | Concentration | Compound                  | Concent | ration |
|--------------------------|---------------|---------------------------|---------|--------|
| acrolein                 | 10 U          | 1,1-dichloroethene        | 5       | U      |
| acrylonitrile            | 10 U          | cis-1,2-dichloroethene    | 5       | U      |
| benzene                  | 72            | trans-1,2-dichloroethene  | 5       | Ù      |
| bromodichloromethane     | 3 · ¹ J       | 1,2-dichloropropane       | 5       | U      |
| bromoform                | 5 U           | cis-1,3-dichloropropene   | 5       | Ű      |
| bromomethane             | 10 U          | trans-1,3-dichloropropene | 5       | U      |
| carbon tetrachloride     | 5 U           | ethyl benzene             | 30      |        |
| chlorobenzene            | 5 U           | methylene chloride        | 4,700   | DB     |
| chloroethane             | 10 U          | 1,1,2,2-tetrachloroethane | . 5     | U      |
| 2-chloroethylvinyl ether | 10 U          | tetrachloroethene         | 5       | U      |
| chloroform               | 62            | toluene                   | 240     | D      |
| chloromethane            | 10 Ü          | 1,1,1-trichloroethane     | 2       | J      |
| dibromochloromethane     | 5 U           | 1,1,2-trichloroethane     | 5       | U      |
| 1,1-dichloroethane       | 5 U           | trichloroethene           | 5       | U      |
| 1,2-dichloroethane       | 5 Û           | trichlorofluoromethane    | 5       | U      |
|                          |               | vinyl chloride            | 10      | U      |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/19/92

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

D - Compound analyzed at a secondary dilution factor.

Client Project ID: LE Carpenter Job Number: ITPK 51499

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in Total µg

Sample Matrix: VOST Tube

Client Sample ID: 1045-069-07T Lab Sample ID: K3272T

| <u>Tentative Identification (1)</u> | Concentration (2) |  |  |
|-------------------------------------|-------------------|--|--|
| ethane, 1,1,2-trichloro-1,2,        | 15                |  |  |
| benzene, dimethyl-                  | 900 Y             |  |  |
| benzene, 1,2-dimethyl-              | 120               |  |  |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.
- Y Indistinguishable isomer in tentatively identified compounds.

Job Number: ITPK 51499

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in Total  $\mu g$ Sample Matrix: VOST Tube

Client Sample ID: 1045-069-07M Lab Sample ID: K3272M

| Compound                 | Concentration |   | Compound                  | Concent | Concentration |  |
|--------------------------|---------------|---|---------------------------|---------|---------------|--|
| acrolein                 | 10            | Ü | 1,1-dichloroethene        | 5       | U             |  |
| acrylonitrile            | 10            | U | cis-1,2-dichloroethene    | . 5     | Ü             |  |
| benzene                  | 5             | U | trans-1,2-dichloroethene  | 5       | U             |  |
| bromodichloromethane     | 5             | U | 1,2-dichloropropane       | 5       | Ü             |  |
| bromoform                | 5             | U | cis-1,3-dichloropropene   | . 5     | U ·           |  |
| bromomethane             | 10            | U | trans-1,3-dichloropropene | 5       | U             |  |
| carbon tetrachloride     | 5             | U | ethyl benzene             | 5       | U             |  |
| chlorobenzene            | 5             | U | methylene chloride        | 4,600   | DB            |  |
| chloroethane             | 10            | Ų | 1,1,2,2-tetrachloroethane | 5       | U             |  |
| 2-chloroethylvinyl ether | 10            | U | tetrachloroethene         | 5       | U             |  |
| chloroform               | 6             |   | toluene                   | 5       | U             |  |
| chloromethane            | 10            | U | 1,1,1-trichloroethane     | 5       | Ü             |  |
| dibromochloromethane     | 5             | U | 1,1,2-trichloroethane     | 5       | Ū             |  |
| 1,1-dichloroethane       | 5             | U | trichloroethene           | . 5     | U             |  |
| 1,2-dichloroethane       | 5             | U | trichlorofluoromethane    | 5       | Ü             |  |
|                          |               |   | vinyl chloride            | 10      | Ŭ             |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/22/92

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

D - Compound analyzed at a secondary dilution factor.

Client Project ID: LE Carpenter

Job Number: ITPK 51499

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in Total µg

Sample Matrix: VOST Tube

Client Sample ID: 1045-069-07M Lab Sample ID: K3272M

| Tentative Identification (1) | Concentration (2) |
|------------------------------|-------------------|
| ethane, 1,1,2-trichloro-1,2, | 5.5               |
| carbon disulfide             | 6.7               |

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Client Project ID: LE Carpenter

Job Number: ITPK 51499

### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in Total  $\mu g$ Sample Matrix: VOST Tube

Client Sample ID: 1045-069-08T Lab Sample ID: K3273T

| Compound                 | Concentration | Compound                  | Concentration     |  |
|--------------------------|---------------|---------------------------|-------------------|--|
| acrolein                 | 10 U          | 1,1-dichloroethene        | 5 U               |  |
| acrylonitrile            | 10 U          | cis-1,2-dichloroethene    | 5 U               |  |
| benzene                  | <b>5</b> Ŭ    | trans-1,2-dichloroethene  | 5 U               |  |
| bromodichloromethane     | 5 U           | 1,2-dichloropropane       | 5 U               |  |
| bromoform                | 5 U           | cis-1,3-dichloropropene   | <i>5</i> <b>U</b> |  |
| bromomethane             | 10 U          | trans-1,3-dichloropropene | 5 U               |  |
| carbon tetrachloride     | , <b>5</b> U  | ethyl benzene             | 5 U               |  |
| chlorobenzene            | 5 : Ŭ         | methylene chloride        | 5 B               |  |
| chloroethane             | 10 U          | 1,1,2,2-tetrachloroethane | 5 U               |  |
| 2-chloroethylvinyl ether | 10 U          | tetrachloroethene         | 5 U               |  |
| chloroform               | 5 Ų           | toluene                   | 1 J               |  |
| chloromethane            | 10 U          | 1,1,1-trichloroethane     | 5 U               |  |
| dibromochloromethane     | 5 (U          | 1,1,2-trichloroethane     | 5 U               |  |
| 1,1-dichloroethane       | 5 U           | trichloroethene           | 5 U               |  |
| 1,2-dichloroethane       | 5 U           | trichlorofluoromethane    | 5 U               |  |
|                          | e* *          | vinyl chloride            | 10 · U            |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/19/92

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

Job Number: ITPK 51499

Client Project ID: LE Carpenter

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in Total µg

Sample Matrix: VOST Tube

Client Sample ID: 1045-069-08T Lab Sample ID: K3273T

Tentative Identification (1)

Concentration (2)

None Detected

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITPK 51499

#### **VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS**

Results in Total  $\mu g$ Sample Matrix: VOST Tube

Client Sample ID: 1045-069-08M Lab Sample ID: K3273M

| Compound                 | Concentration | Compound                  | Concentration |
|--------------------------|---------------|---------------------------|---------------|
| acrolein                 | 10 U          | 1,1-dichloroethene        | 5 U           |
| acrylonitrile            | 10 U          | cis-1,2-dichloroethene    | 5 U           |
| benzene                  | 5 U           | trans-1,2-dichloroethene  | 5 U           |
| bromodichloromethane     | 5 . U         | 1,2-dichloropropane       | 5 U           |
| bromoform                | 5 U           | cis-1,3-dichloropropene   | 5 U           |
| bromomethane             | 10 U          | trans-1,3-dichloropropene | 5 U           |
| carbon tetrachloride     | 5 . U         | ethyl benzené             | 5 U           |
| chlorobenzene            | 5 U           | methylene chloride        | 8 B           |
| chloroethane             | 10 U          | 1,1,2,2-tetrachloroethane | 5 U           |
| 2-chloroethylvinyl ether | 10 U          | tetrachloroethene         | 5 U           |
| chloroform               | 5 U           | toluene                   | 5 U           |
| chloromethane            | 10 U          | 1,1,1-trichloroethane     | 5 U           |
| dibromochloromethane     | <b>5</b> L Ü  | 1,1,2-trichloroethane     | 5 U           |
| 1,1-dichloroethane       | 5 U           | trichloroethene           | 5 U           |
| 1,2-dichloroethane       | 5 U           | trichlorofluoromethane    | 5 U           |
|                          |               | vinyl chloride            | 10 U          |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/22/92

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

IT ANALYTICAL SERVICES 5815 MIDDLEBROOK PIKE KNOXVILLE, TN

Job Number: ITPK 51499

#### ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in Total µg

Sample Matrix: VOST Tube

Client Sample ID: 1045-069-08M Lab Sample ID: K3273M

Tentative Identification (1)

Concentration (2)

None Detected

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Job Number: ITPK 51499

# VOLATILE ORGANIC PRIORITY POLLUTANT ANALYSIS Results in Total µg

Sample Matrix: VOST Tube

Client Sample ID: 1045-069-08B Lab Sample ID: K3273B

| Compound                 | Concentration |   | Compound                  | Concent | Concentration |  |
|--------------------------|---------------|---|---------------------------|---------|---------------|--|
| acrolein                 | 10            | U | 1,1-dichloroethene        | 5       | U             |  |
| acrylonitrile            | 10            | U | cis-1,2-dichloroethene    | 5       | U             |  |
| benzene                  | 6             |   | trans-1,2-dichloroethene  | 5       | U             |  |
| bromodichloromethane     | 5             | U | 1,2-dichloropropane       | 5       | U             |  |
| bromoform                | 5             | U | cis-1,3-dichloropropene   | 5       | U             |  |
| bromomethane             | 10            | Ü | trans-1,3-dichloropropene | 5       | U             |  |
| carbon tetrachloride     | 5             | U | ethyl benzene             | 10      |               |  |
| chlorobenzene            | 5             | U | methylene chloride        | 20      | В             |  |
| chloroethane             | 10            | U | 1,1,2,2-tetrachloroethane | 5       | U             |  |
| 2-chloroethylvinyl ether | 10            | U | tetrachloroethene         | . 5     | U.            |  |
| chloroform               | 11            |   | toluene                   | 10      |               |  |
| chloromethane            | .10           | U | 1,1,1-trichloroethane     | 5       | U             |  |
| dibromochloromethane     | 5             | U | 1,1,2-trichloroethane     | 5       | U             |  |
| 1,1-dichloroethane       | .5            | U | trichloroethene           | 5       | U             |  |
| 1,2-dichloroethane       | 5             | U | trichlorofluoromethane    | 5       | U             |  |
|                          |               | 1 | vinyl chloride            | 10      | U             |  |

U - Compound was analyzed for but not detected. The number is the detection limit for the sample.

Date of Analysis: 06/23/92

J - Indicates an estimated value less than the detection limit.

B - Analyte was found in the blank as well as the sample.

KNOXVILLE, TN

Job Number: ITPK 51499

## ADDITIONAL VOLATILE ORGANIC COMPOUNDS

Results in Total µg

Sample Matrix: VOST Tube

Client Sample ID: 1045-069-08B Lab Sample ID: K3273B

Client Project ID: LE Carpenter

| Tentative Identification (1) | Concentration (2) |
|------------------------------|-------------------|
| pentane                      | 2,300             |
| hexane                       | 1,300             |
| octane                       | 290               |
| benzene, dimethyl-           | 1,600 Y           |
| benzene, 1,2-dimethyl-       | 320               |

Remarks:

- (1) Identification is based on computer search of the NIST Library.
- (2) Concentration is based on a response factor of 1.00 relative to the internal standard.

Y - Indistinguishable isomer in tentatively identified compounds.



### CERTIFICATE OF ANALYSIS

**IT** Corporation

312 Directors Drive

Knoxville, Tennessee 37923

Attn: Michael Krstich

Job Number: ITAD 51019

P.O. Number: 483048

May 1, 1992

F.O. Mullioet. 403040

This is the Certificate of Analysis for the following samples:

Client Project ID:

LE Carpenter

Date Received by Lab:

03/26/92

Number of Samples:

Eight (8)

Sample Type:

Water - five (5), Soil - three (3)

#### I. Introduction

On 03/26/92, five (5) water samples and three (3) soil samples arrived at the ITAS-Knoxville, Tennessee, laboratory from the IT-Technology Development Laboratory, Knoxville, Tennessee, in support of the LE Carpenter project. The list of analytical tests performed, as well as date of receipt and analysis, can be found in the attached report.

#### II. Analytical Results/Methodology

The analytical results for this report are presented by analytical test. Each set of data will include sample identification information and the analytical results. Please note that the data are not blank corrected.

The samples were analyzed for priority pollutant volatile organic compounds by gas chromatography/mass spectroscopy (GC/MS) based on EPA SW-846 method 8240.

The samples were analyzed for priority pollutant semivolatile organic compounds by gas chromatography/mass spectroscopy (GC/MS) based on EPA SW-846 method 8270.

Reviewed and Approved:

Alyce R. Moore

Client Project ID: LE Carpenter

Job Number: ITAD 51019

#### II. Analytical Results/Methodology (continued)

Chemical oxygen demand (COD) was measured using the EPA approved HACH procedure, <u>HACH Water Analysis Handbook</u>, HACH Chemical Company, 1980.

#### III. Quality Control

Routine laboratory level I QC was followed.

The volatiles analyses were performed by purge and trap with a J&W DB-624 megabore column on a Finnigan OWA GC/MS/DS. The sample runs generally went well; however, samples 1045-055-2 and 1045-055-3 required necessary dilutions for methylene chloride and toluene. Sample 1045-052-1 had low toluene-d8 and dichloroethane-d4 surrogate recoveries. This appeared to have been due to matrix interference since a reanalysis showed similar results and additional samples analyzed after these samples had acceptable surrogate recoveries. Xylene was not a target and was reported as a TIC, but was quantified similar to a target compound using a continuing calibration standard. Sample 1045-055-1 required a dilution for xylene.

The semivolatiles analyses were performed by direct injection of sample extract on a Restek XTI-5 capillary column on a Finnigan 4500 GC/MS/DS. The sample runs generally went well; however, high levels of various target phthalates were seen requiring dilutions. Two dilutions were required to quantify both bis(2-ethylhexyl)phthalate and butylbenzyl phthalate in sample 1045-052-3. The bis(2-ethylhexyl) phthalate calibration curve was slightly exceeded in the dilution of sample 1045-052-1. This parameter has, in our experience, been linear beyond the calibration range on this instrument, and the data was judged as reliable. High levels of xylene were seen in the BN extracts. There were no problems seen in final data review.

The COD of the samples was measured using the reactor digestion method followed by titration with ferrous ammonium sulfate. No problems were encountered.



## CERTIFICATE OF ANALYSIS

IT Corporation

312 Directors Drive

Knoxville, Tennessee 37923

Attn: Michael Krstich

Job Number: ITAD 51191

P.O. Number:

May 1, 1992

This is the Certificate of Analysis for the following samples:

Client Project ID:

LE Carpenter

1

Date Received by Lab:

04/14/92

Number of Samples:

Six (6)

Sample Type:

Soil

#### I. Introduction

On 04/14/92, six (6) soil samples arrived at the ITAS-Knoxville, Tennessee, laboratory from the IT-Technology Development Laboratory, Knoxville, Tennessee, in support of the LE Carpenter project. The list of analytical tests performed, as well as date of receipt and analysis, can be found in the attached report.

#### II. Analytical Results/Methodology

The analytical results for this report are presented by analytical test. Each set of data will include sample identification information and the analytical results. Please note that the data are not blank corrected.

The samples were analyzed for total organic carbon (TOC) using modified EPA method 9060. The results should be considered nonpurgeable organic carbon since the sample preparation process may result in loss of volatiles.

Reviewed and Approved:

Alyce R. Moore

Client Project ID: LE Carpenter

Job Number: ITAD 51191

#### III. Quality Control

Routine laboratory level I QC was followed.

The TOC content of the samples was determined by chemical wet oxidation followed by infrared detection. Samples were weighed into glass ampules and mixed with a phosphoric acid/sodium persulfate solution. The phosphoric acid converts all inorganic carbon to carbon dioxide (CO<sub>2</sub>) which is purged from the ampule with oxygen. The ampule is then sealed under oxygen and heated to approximately 100°C in a water bath for 30 minutes. The heating process converts the organic carbon to CO<sub>2</sub> which is trapped in the sealed ampule. The samples were analyzed by an O.I. Corporation total organic carbon analyzer, Model 700; the carbon was quantified using infrared detection by measuring the absorbance of CO<sub>2</sub>. No major problems were encountered.



### CERTIFICATE OF ANALYSIS

IT Corporation

312 Directors Drive

Knoxville, Tennessee 37923

Attn: Michael Krstich

Job Number: ITAD 51254

May 29, 1992

P.O. Number: 483100,001

This is the Certificate of Analysis for the following samples:

Client Project ID:

LE Carpenter

Date Received by Lab:

04/24/92

Number of Samples:

Four (4)

Sample Type:

Water

#### I. Introduction

On 04/24/92, four (4) water samples arrived at the ITAS-Knoxville, Tennessee, laboratory from the IT-Technology Development Laboratory, Knoxville, Tennessee, in support of the LE Carpenter project. The list of analytical tests performed, as well as date of receipt and analysis, can be found in the attached report.

#### II. Analytical Results/Methodology

The analytical results for this report are presented by analytical test. Each set of data will include sample identification information and the analytical results. Please note that the data are not blank corrected.

The samples were analyzed for priority pollutant volatile organic compounds by gas chromatography/mass spectroscopy (GC/MS) based on EPA SW-846 method 8240.

The samples were analyzed for priority pollutant semivolatile organic compounds by gas chromatography/mass spectroscopy (GC/MS) based on EPA SW-846 method 8270.

Reviewed and Approved:

Alvce R. Moore

Laboratory Manager

American Council of Independent Laboratories International Association of Environmental Testing Laboratories American Association for Laboratory Accreditation

Client Project ID: LE Carpenter

Job Number: ITAD 51254

#### II. Analytical Results/Methodology (continued)

Chemical oxygen demand (COD) was measured using the EPA approved HACH procedure, <u>HACH Water Analysis Handbook</u>, HACH Chemical Company, 1980.

#### III. Quality Control

Routine laboratory level I QC was followed.

The volatiles analyses were performed by purge and trap with a J&W DB-624 megabore column on a Finnigan OWA GC/MS/DS. The sample runs went well. Methylene chloride was found in several samples at levels requiring dilutions. Bromofluorobenzene was slightly high in undiluted analysis of sample 1045-061-1, but was compliant in the dilution, indicating a slight sample matrix effect. Overall analysis and method QC looked good. The semivolatiles analyses were performed by direct injection of sample extract on a Restek XTI-5 capillary column on a Fisons VG TRIO-1 GC/MS/DS. The sample runs went well. There were no problems seen in final data review.

The COD of the samples was measured using the reactor digestion method followed by titration with ferrous ammonium sulfate. No major problems were encountered.



### CERTIFICATE OF ANALYSIS

IT Corporation

312 Directors Drive

Knoxville, Tennessee 37923

Attn: Michael Krstich

Job Number: ITAD 51315

May 27, 1992

P.O. Number: 483100.001

This is the Certificate of Analysis for the following samples:

Client Project ID:

LE Carpenter

Date Received by Lab:

05/05/92

Number of Samples:

Four (4)

Sample Type:

Water

#### I. Introduction

On 05/05/92, four (4) water samples arrived at the ITAS-Knoxville, Tennessee, laboratory from the IT-Technology Development Laboratory, Knoxville, Tennessee, in support of the LE Carpenter project. The list of analytical tests performed, as well as date of receipt and analysis, can be found in the attached report.

#### II. Analytical Results/Methodology

The analytical results for this report are presented by analytical test. Each set of data will include sample identification information and the analytical results. Please note that the data are not blank corrected.

The samples were analyzed for priority pollutant volatile organic compounds by gas chromatography/mass spectroscopy (GC/MS) based on EPA SW-846 method 8240.

The samples were analyzed for priority pollutant semivolatile organic compounds by gas chromatography/mass spectroscopy (GC/MS) based on EPA SW-846 method 8270.

Reviewed and Approved:

Alyce R. Moore

Laboratory Manager

American Council of Independent Laboratories International Association of Environmental Testing Laboratories American Association for Laboratory Accreditation

Client Project ID: LE Carpenter

Job Number: ITAD 51315

#### II. Analytical Results/Methodology (continued)

Chemical oxygen demand (COD) was measured using the EPA approved HACH procedure, <u>HACH Water Analysis Handbook</u>, HACH Chemical Company, 1980.

#### III. Quality Control

Routine laboratory level I QC was followed.

The volatiles analyses were performed by purge and trap with a J&W DB-624 megabore column on two Finnigan OWA GC/MS/DS units. The sample runs generally went well, however, interferences such as organic contaminants and surfactants did cause some problems. More concentrated analyses than reported were performed, but the interferences encountered rendered the data unreliable. For example, sample 1045-063-1 required a 100-fold dilution for successful analysis due to high surfactant levels and background from substituted benzenes and other hydrocarbons. Some dilutions were also required for quantification, e.g., sample 1045-063-3 was diluted 25-fold for methylene chloride. Overall, the method QC and analyses looked good.

The semivolatiles analyses were performed by direct injection of sample extract on a Restek XTI-5 capillary column on a Fisons VG TRIO-1 GC/MS/DS. The sample runs generally went well, however, some problems were encountered due to high levels of organic contamination in the samples. Originally a 1,000 ml aliquot of each sample was extracted, but sample 1045-063-1 and 1045-063-4 would not concentrate to the method required volume of 1.0 ml. Further dilutions would be required for analysis and surrogate data for these two samples would be lost, hence, reextractions of these samples were performed using 20 ml aliquots resulting in 50-fold dilutions with measurable surrogate recoveries. A further fivefold dilution was required of sample 1045-063-1 to quantify bis(2-ethylhexyl) phthalate. The high level of bis(2-ethylhexyl) phthalate in the original analysis of 1045-063-1 suppressed one of the internal standards resulting in one low internal standard area recovery and one high surrogate recovery. These anomalies disappeared in the subsequent dilution. There were no problems seen in final data review.

The COD of the samples was measured using the reactor digestion method followed by titration with ferrous ammonium sulfate. No major problems were encountered.



## CERTIFICATE OF ANALYSIS

IT Corporation

312 Directors Drive

Knoxville, Tennessee 37923

Attn: Michael Krstich

Job Number: ITAD 51328

May 26, 1992

P.O. Number: NA

This is the Certificate of Analysis for the following sample:

Client Project ID:

LE Carpenter

Date Received by Lab:

05/06/92

Number of Samples:

One (1)

Sample Type:

Soil

#### I. Introduction

On 05/06/92, one (1) soil sample arrived at the ITAS-Knoxville, Tennessee, laboratory from the ITrechnology Development Laboratory, Knoxville, Tennessee, in support of the LE Carpenter project. The list of analytical tests performed, as well as date of receipt and analysis, can be found in the attached report.

#### II. Analytical Results/Methodology

The analytical results for this report are presented by analytical test. Each set of data will include sample identification information and the analytical results. Please note that the data are not blank corrected.

The sample was analyzed for total organic carbon (TOC) using modified EPA method 9060. The results should be considered nonpurgeable organic carbon since the sample preparation process may result in loss of volatiles.

Reviewed and Approved:

Alyce R. Moore

Client Project ID: LE Carpenter

Job Number: ITAD 51328

#### III. Quality Control

Routine laboratory level I QC was followed.

The TOC content of the sample was determined by chemical wet oxidation followed by infrared detection. The sample was weighed into a glass ampule and mixed with a phosphoric acid/sodium persulfate solution. The phosphoric acid converts all inorganic carbon to carbon dioxide (CO2) which is purged from the ampule with oxygen. The ampule is then sealed under oxygen and heated to approximately 100°C in a water bath for 30 minutes. The heating process converts the organic carbon to CO2 which is trapped in the sealed ampule. The sample was analyzed by an O.I. Corporation total organic carbon analyzer, Model 700; the carbon was quantified using infrared detection by measuring the absorbance of CO2. No major problems were encountered.



### CERTIFICATE OF ANALYSIS

IT Corporation

June 23, 1992

312 Directors Drive Knoxville, TN 37923 Attn: Michael Krstich

Job Number: ITPK 51491

P.O. Number: 483100.001

This is the Certificate of Analysis for the following samples:

Client Project ID:

LE Carpenter

Date Received by Lab:

06/01/92

Number of Samples:

Twelve (12)

Sample Type:

Water - eight (8), Soil - four (4)

#### I. Introduction

On 06/01/92, eight (8) water samples and four (4) soil samples arrived at the ITAS-Knoxville, Tennessee, laboratory from IT Corporation, Knoxville, Tennessee, in support of the LE Carpenter project. The list of analytical tests performed, as well as date of receipt and analysis, can be found in the attached report.

#### II. Analytical Results/Methodology

The analytical results for this report are presented by analytical test. Each set of data will include sample identification information and the analytical results. Please note that the data are not blank corrected, and soil samples were reported on a dry weight basis.

The samples were analyzed for priority pollutant volatile organic compounds by gas chromatography/mass spectroscopy (GC/MS) based on EPA SW-846 method 8240.

The samples were analyzed for priority pollutant semivolatile organic compounds by gas chromatography/mass spectroscopy (GC/MS) based on EPA SW-846 method 8270.

Reviewed and Approved:

Robert L. Curry

Laboratory Systems Manager

American Council of Independent Laboratories
International Association of Environmental Testing Laboratories
American Association for Laboratory Accreditation

Client Project ID: LE Carpenter

Job Number: ITPK 51491

#### II. Analytical Results/Methodology (continued)

The samples were analyzed for total organic carbon (TOC) based on EPA methods 415.1 and 9060.

Chemical oxygen demand (COD) was measured using the EPA approved HACH procedure, <u>HACH Water Analysis Handbook</u>, HACH Chemical Company, 1980.

#### III. Quality Control

Routine laboratory level I QC was followed.

The volatiles analyses were performed by purge and trap with a J&W DB-624 megabore column on a Finnigan INCOS 500 and OWA GC/MS/DS units. The sample runs went well, although there were several additional dilutions required. The initial run of 1045-068-01 showed one surrogate marginally outside limits, while the dilution met limits. The TIC analysis showed primarily cyclics, alkanes, and alkylbenzenes; the quality of the matches was better in some runs than others. The matches were tentative, and listed names should be taken as guides to compound type.

The semivolatiles analyses were performed by direct injection of sample extract on a Restek XTI-5 capillary column on a VG TRIO-1 GC/MS/DS. The sample runs went well with the exception of the original run of 10456-067-01, in which some internal standard and surrogate variants were seen. The deviations were probably due to matrix effect from the high levels in the sample. Many extracts required dilution prior to analysis. As the samples contained a large number of components, TIC analysis was fairly involved: as for volatiles, it is suggested that the listed names be taken as guides to compound type, as the peaks sometimes had more than one reasonable spectral match, and also some spectra may have been overlapped in the generally crowded chromatogram. A number of phthalates were seen; in particular, bis(2-ethylhexyl) eluted in a general phthalate cluster, and judgement was used to distinguish the priority pollutant from the TICs (please note the TIC phthalates around 28.3 minutes which all showed similar spectra to the target compound, but were judged to be from similar but separate compounds). There were no other problems seen in final data review.



## CERTIFICATE OF ANALYSIS

IT Corporation

312 Directors Drive

Knoxville, TN 37923

Attn: Michael Krstich

Job Number: ITPK 51499

June 30, 1992

P.O. Number: 483100.001

This is the Certificate of Analysis for the following samples:

Client Project ID:

LE Carpenter

Date Received by Lab:

06/02/92

Number of Samples:

Eight (8)

Sample Type:

Water - four (4), XAD Tubes - four (4)

#### I. Introduction

On 06/02/92, four (4) water samples and four (4) XAD tubes arrived at the ITAS-Knoxville, Tennessee, laboratory from IT Corporation, Knoxville, Tennessee, in support of the LE Carpenter project. The list of analytical tests performed, as well as date of receipt and analysis, can be found in the attached report.

#### II. Analytical Results/Methodology

The analytical results for this report are presented by analytical test. Each set of data will include sample identification information and the analytical results. Please note that the data are not blank corrected.

The samples were analyzed for priority pollutant volatile organic compounds by gas chromatography/mass spectroscopy (GC/MS) based on EPA SW-846 method 8240.

The water samples were analyzed for priority pollutant semivolatile organic compounds by gas chromatography/mass spectroscopy (GC/MS) based on EPA SW-846 method 8270.

Reviewed and Approved:

Alyce R. Moore

Job Number: ITPK 51499

Client Project ID: LE Carpenter

#### III. Quality Control

Routine laboratory level I QC was followed.

The volatiles analyses on VOST tubes were performed by purge and trap with a J&W DB-624 megabore column on a Finnigan OWA GC/MS/DS. The sample runs generally went well considering the charcoal in the matrix. As expected, the surrogate recoveries were low compared to normal recovery ranges. Three VOST tubes were received for each sample: 1045-069-05, 1045-069-06 and 1045-069-08 labeled T, M, and. Two VOST tubes were received for 1045-069-07 labeled T and M. A single letter suffix was added to the sample names corresponding to the T, M, or B label. The B tubes exhibited the highest levels of target compounds requiring dilutions in most cases. Sample 1045-069-05B not only showed poor surrogate recovery, but also yielded a low internal standard area. This phenomenon was not present in the more dilute analyses and was considered an effect of matrix. The volatiles analyses on water samples were performed by purge and trap with a J&W DB-624 megabore column on a Finnigan INCOS 500 GC/MS/DS. The sample runs went well; sample 1045-069-01 required dilution because of excessive foaming with the intital run. The TIC analysis was generally straightforward, but it is suggested that the listed names be used as guides to compound type, as there were often (both here and in semivolatiles, below) comparable matches from more than one database entry.

The semivolatiles analyses were performed by direct injection of sample extract on a Restek XTI-5 capillary column on a Finnigan 4500 GC/MS/DS. The sample runs generally went well. A reextraction of sample 1045-069-04 by liquid/liquid extraction was performed due to some low base/neutral surrogate recoveries. The reextraction could only be concentrated to 5.0 ml methylene chloride compared to the normally achievable 1.0 ml due to higher boiling fractions in the sample. Both sets of data were reported for comparison. The TIC analysis showed a number of unusual compounds, many with poor spectral matches against the EPA NIST database. The spectra for many unknowns appeared to be of "O" containing alkanes; of particular note was that many spectra strongly suggested "crown" compounds, although the specific structure and type were uncertain. There were no other problems seen in final data review.